Full Raw Water
Quality Characterization

Step 2 of 97-005 Evaluation

City of Santa Monica

Olympic Well Field

August 2021

308038-03533













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# **Executive Summary**

As groundwater in the vicinity of the Olympic Well Field is impaired by contamination, the City of Santa Monica (the City) is required to demonstrate compliance with the Division of Drinking Water (DDW) Process Memo 97-005-R2020 - *Revised Guidance for Direct Domestic Use of Extremely Impaired Sources*. A source is considered "extremely impaired" if it meets two or more of 10 DDW-developed criteria. Based on the evaluation of available water quality data, groundwater in the vicinity of the Olympic Well Field meets up to three criteria.

The DDW 97-005 evaluation process consists of 10 steps for assessing proposals, establishing appropriate permit conditions, and approving the use of an extremely impaired drinking water source. This report documents the second step (Step 2) of the 97-005 evaluation process, i.e., "Full Characterization of the Raw Water Quality" for the Olympic Well Field. The purpose of this step is the to fully characterize constituents in the raw water produced by the Olympic Well Field, thereby ensuring a treatment system can be properly designed. In accordance with the DDW Process Memo 97-005-R2020, Step 2 (this report) was conducted after finalization of Step 1 (Drinking Water Source Assessment and Contaminant Assessment), as "each step relies upon the findings and conclusions of the prior step" (DDW 2020).

A series of data analyses were conducted to evaluate raw water quality to meet the primary objective of characterizing influent water quality that will enter the planned treatment system so that an appropriate level of monitoring and treatment can be designed. This included screening of water quality data against specific criteria and regulatory values to identify constituents of potential concern (COPCs), analyzing water quality data to estimate future treatment plant influent concentrations, trend analysis for historical water quality to understand how water quality trends have changed over time, and analysis of variability to develop an understanding of how water quality has changed under the influence of certain factors such as pumping and seasonal variation in precipitation.

An initial screening of water quality data indicated 42 COPCs based on constituents detected in production wells (or constituents with three or more detections in monitoring wells), with a ratio of maximum concentration to maximum contaminant level (MCL) or notification level (NL) greater than 0.05 (5%). The list of COPCs was then further refined to determine which would drive treatment system design (using ultraviolet/advanced oxidation process [UV/AOP] and granular activated carbon [GAC]) based on constituents which are synthetic organics and had a ratio of maximum concentration to MCL or NL greater than 0.5 (50%). This yielded a list of 15 synthetic organic COPCs, as follows:

- Nine COPCs from production well data:
  - 1,1-dichloroethane (1,1-DCA)
  - 1,1-dichloroethene (1,1-DCE)
  - 1,2,3-trichloropropane (1,2,3-TCP)
  - 1.4-dioxane
  - carbon tetrachloride
  - cis-1,2-dichloroethene (cis-1,2-DCE)
  - perfluorooctanoic acid (PFOA)
  - tetrachloroethene (PCE)
  - trichloroethene (TCE)

- An additional six COPCs from monitoring well data:
  - 1,1,2-trichloroethane (1,1,2-TCA)
  - 1,2-dichloroethane (1,1-DCA)
  - benzene
  - methyl tert-butyl ether (MTBE)
  - trans-1,2-dichloroethene (trans-1,2-DCE)
  - vinyl chloride







Statistical analysis and flow-weighting calculations were then conducted to estimate future treatment plant influent concentrations using monitoring well groundwater quality data for all constituents including the identified COPCs. The results of this analysis indicated four COPCs are projected to be at concentrations above their respective MCL or NL in treatment plant influent, as summarized in Table E-1. These COPCs are 1,4-dioxane, PCE, TCE and 1,2,3-TCP. For comparison, calculated influent estimates based on recent (2020) water quality data from production wells SM-4, SM-8 and SM-9 indicate that 1,4-dioxane, PCE and TCE in influent would be above their respective MCL or NL, with 1,2,3-TCP being below its MCL based on these data, as shown in Table E-1.

Table E-1 Summary of Estimated Concentrations in Treatment Plant Influent for Key Synthetic Organic COPCs

		Using UC	L95*	Using Production Well Concentrations From 2020**			
Constituent [Units]	MCL or NL	Plant Influent Concentration Estimates	With Safety Factor Applied***	Plant Influent Concentration Estimates	With Safety Factor Applied***		
1,2,3-TCP [μg/L] 0.00		0.022 0.026		0.00091	0.0011		
1,4-Dioxane [µg/L]	1	13.8	20.7	9.9	14.9		
PCE [µg/L]	5	10.4	15.6	31	46.5		
TCE [μg/L]	5	8.2	12.3	23	34.5		

**Notes:**  $\mu g/L = micrograms per liter; values above respective MCLs or NLs are highlighted$ **bold**.

Trend analysis was performed on available temporal production well data for COPCs to inform potential future concentration trends and design of the treatment system. Key outcomes from the trend analysis include the identification of statistically significant increasing trends for PCE and 1,4-dioxane. Although no statistically significant trend was identified for TCE, visual assessment of trend charts indicates that recent (2018-2020) concentrations are rebounding to the elevated levels recorded between 2012 and 2014. No statistically significant or visual trend was identified for 1,2,3-TCP. Visual assessment indicated nitrate and nitrite (as nitrogen) exhibited variability since start of analysis period (2012) with no discernible trend identified. Visual assessment of other COPC trends generally indicated that parameters with concentrations above detection limits remained stable during the analysis period (2012-2020).

An assessment of concentration variability generally indicates that concentrations of 1,4-dioxane, TCE, PCE exhibit an observable correlation to operational status (i.e., pumping or non-pumping) and pumping rate of a well or at nearby wells; however, seasonal impacts or seasonal signals are not distinguishable in the data sets.

<sup>\*</sup> From production wells concentration estimates.

<sup>\*\*</sup>Maximum observed values from production wells from available 2020 sampling data (provided for information only).

<sup>\*\*\*</sup>A safety factor of 1.5 was applied for each constituent, except for 1,2,3-TCP which used a safety factor of 1.2.







# Acronyms and Abbreviations

Acronym / Abbreviation	Definition
%	Percent
1,1,1-TCA	1,1,1-Trichloroethane
1,1-DCA	1,1-Dichloroethane
1,2-DCA	1,2-Dichloroethane
1,1-DCE	1,1-Dichloroethene
1,2,3-TCP	1,2,3-Trichloropropane
μg	Microgram
μg/L	Microgram(s) Per Liter
AST	Aboveground Storage Tank
bgs	Below Ground Surface
CA	Contaminant Assessment
CEPRD	Coalition for Environmental Protection Restoration and Development
cfs	Cubic Foot/Feet Per Second
City	City of Santa Monica
cis-1,2-DCE	cis-1,2-Dichloroethene
COPC	Constituent of Potential Concern
Cr(VI)	Hexavalent Chromium
DDW	Division of Drinking Water
DEHP	Bis(2-ethylhexyl)phthalate
DQO	Data Quality Objective
EPA	United States Environmental Protection Agency
ft	Feet
GAC	Granular Activated Carbon.
GIS	Global Information System
gpm	gallons per minute
Cr(VI)	Hexavalent Chromium
GTE	General Telephone & Electronics Corporation







Acronym / Abbreviation	Definition
HPC	Heterotrophic Plate Count
KM	Kaplan Meier
L	Liter(s)
LA	Los Angeles
MCL	California Maximum Contaminant Level
mg/L	Milligram(s) Per Liter
MRL	Method Reporting Limit
N	Nitrogen
NDMA	N-Nitrosodimethylamine
NL	Notification Level
OCPs	Organochlorine Pesticides
PBCs	Polychlorinated Biphenyls
PCE	Tetrachloroethene
PHG	Public Health Goal
RL	Reporting Limit
RWQC	Raw Water Quality Characterization
S	Statistic
SA	(Drinking Water) Source Assessment
SMCL	California Secondary Maximum Contaminant Level
SOCs	Non-Volatile Synthetic Organic Compounds
SVOCs	Semi-Volatile Organic Compounds
SWRCB	State Water Resources Control Board
TCE	Trichloroethene
TDS	Total Dissolved Solids
TIC	Tentatively Identified Compound
trans-1,2-DCE	trans-1,2-Dichloroethene
UCL95	95 Percent (%) Upper Confidence Limit
USGS	United States Geological Survey
UST	Underground Storage Tank







Acronym / Abbreviation	Definition
UV/AOP	Ultraviolet/Advanced Oxidation Process
VC	Vinyl Chloride
VOCs	Volatile Organic Compounds







# 1 Introduction

The Division of Drinking Water (DDW) 97-005 evaluation process consists of 10 steps (also referred to as elements) for assessing proposals, establishing appropriate permit conditions, and approving the use of an extremely impaired drinking water source. A schematic diagram presenting the 10-step 97-005 evaluation process is presented in Figure 1-1. This report documents the second step (Step 2) of the 97-005 evaluation process, i.e., "Full Characterization of the Raw Water Quality" for the City of Santa Monica's (the City's) Olympic Well Field. For the purposes of this raw water quality characterization (RWQC) study, "raw water" is defined as the groundwater extracted from the Olympic Well Field via three production wells (SM-4, SM-8 and SM-9) which are planned to supply the treatment system.

The Olympic Well Field is one of the City's production well fields situated within the Santa Monica Subbasin. The location of the well field is shown in Figure 1-2. As groundwater in the vicinity of the well field is impaired by contamination, the City is required to demonstrate compliance with the DDW Process Memo 97-005-R2020 - *Revised Guidance for Direct Domestic Use of Extremely Impaired Sources* (DDW 2020). Per the guidance, a source is considered to be "extremely impaired" if it meets two or more of 10 DDW-developed criteria. Based on the evaluation of available water quality data, groundwater in the vicinity of the Olympic Well Field meets up to the following three criteria and is therefore considered extremely impaired:

- Contains a contaminant that exceeds 10 times its notification limit (NL) based on chronic health effects, i.e., 1,4-dioxane.
- Is extremely threatened with contamination due to known contaminating activities within the long term, steady state capture zone of a drinking water well or within the watershed of a surface water intake. As identified in the Drinking Water Source Assessment and Contaminant Assessment (SA/CA; Step 1 of the 97-005 Evaluation; ICF 2020), former Gillette and Boeing facilities, which are located within the Olympic Well Field Study Area, are identified contamination sites.
- Might be considered to contain a mixture of contaminants of health concern beyond what is typically seen in terms of number and concentration of contaminants, i.e., 1,4-dioxane, trichloroethene (TCE), and tetrachloroethene (PCE).

It is imperative that the precursor Step 1 report (SA/CA, ICF 2020) for the Olympic Well Field 97-005 Evaluation is read prior to, or in conjunction with this report, "as each step lies upon the findings and conclusions of the prior step" (DDW 2020).

# 1.1 Olympic Well Field History

Historically, the Olympic Well Field consisted of two active drinking water production wells SM-3 and SM-4 which were installed in 1969 and 1981, respectively. A third production well SM-7, which was never brought on-line, was destroyed and abandoned in March 2018 due to excessive sand production and potential for acting as conduit for contamination migration from the heavily impacted B-Zone to deeper and less impacted aquifer zones. A new production well, SM-8, was drilled and installed in November 2017. More recently, the City replaced SM-3 with a new well, SM-9. The City plans to destroy SM-3 in the near future. Both SM-8 and SM-9 will be brought on-line upon successful amendment of the City's

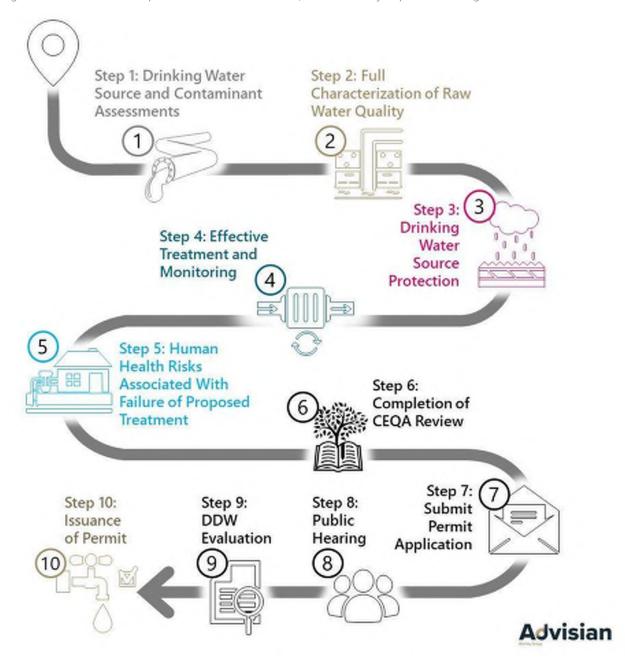


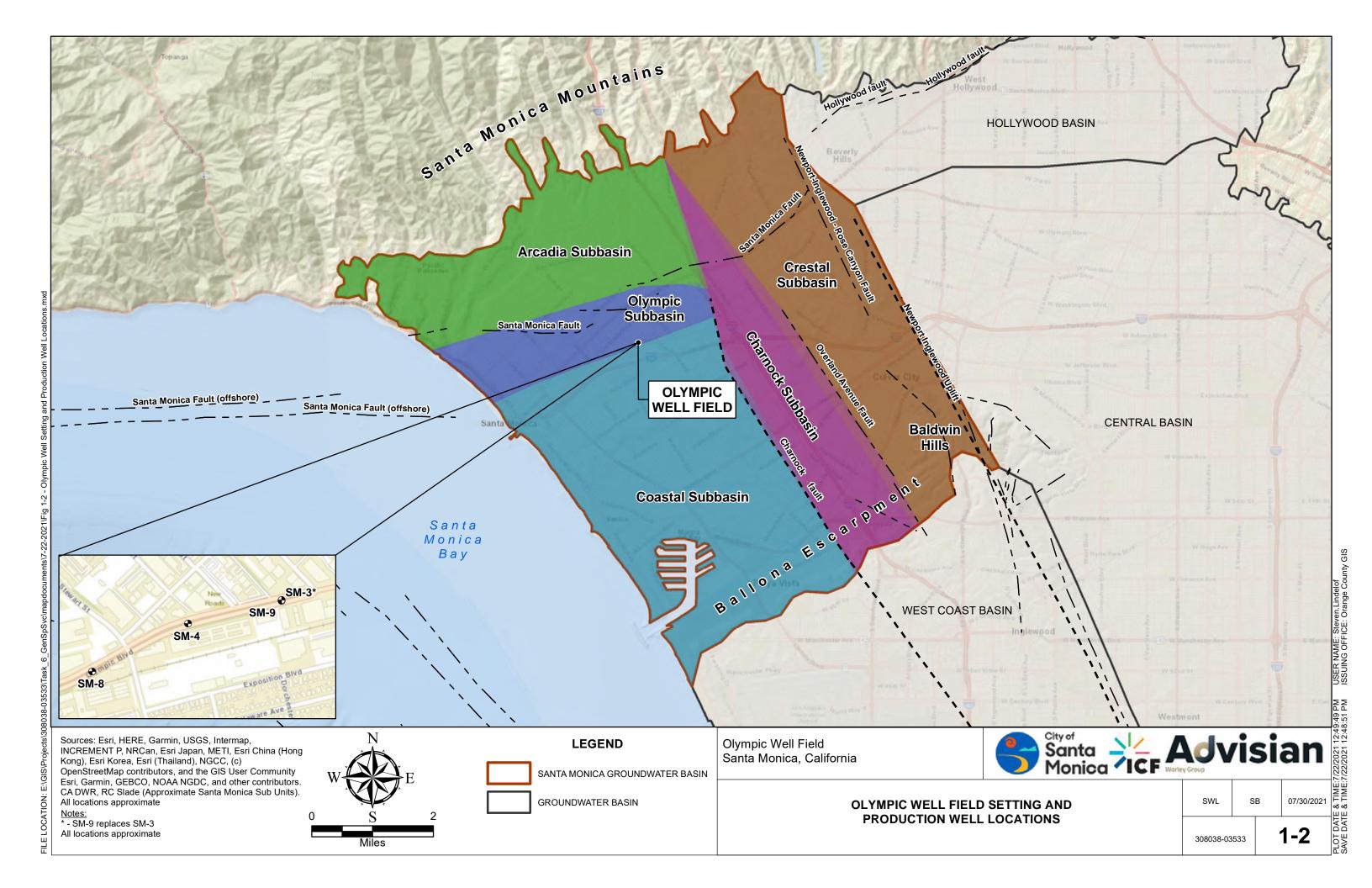




Domestic Water Supply Permit (DDW 1910146). The well field setting and the locations of production wells are shown in Figure 1-2.

Figure 1-1 DDW Ten-Step 97-005 Evaluation Process for an Extremely Impaired Drinking Water Source











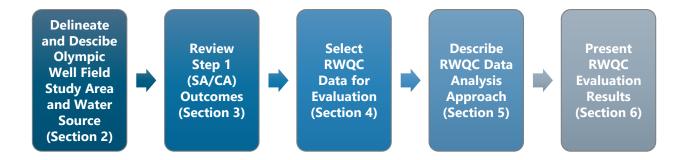
## 1.2 Purpose

The purpose of the RWQC study is to fully characterize constituents in the raw water produced by the Olympic Well Field, thereby ensuring a treatment system can be properly designed. As stated by DDW: "the appropriate level of monitoring and treatment to produce safe drinking water cannot be determined unless the raw water quality is fully understood" (DDW 2020). The outcome of this RWQC study will support the next steps of the 97-005 evaluation process, including "Effective Treatment and Monitoring" (Step 4), which comprises the evaluation and determination of treated water goals and ultimately directs treatment design.

### 1.3 Scope of Work

The scope of work for the RWQC study of the Olympic Well Field is illustrated in Figure 1-3. The Olympic Well Field RWQC study has been conducted in accordance with the DDW Process Memo 97-005-R2020 - Revised Guidance for Direct Domestic Use of Extremely Impaired Sources (DDW 2020).

Figure 1-3 Key Raw Water Quality Characterization Study Tasks



# 1.4 Document Organization

The report is organized into the following sections:

**Section 1 - Introduction:** This section provides the introductory and background information, purpose of study, regulatory requirements and organization of the document.

**Section 2 - Olympic Well Field Water Source:** This section summarizes the characteristics of the Olympic Well Field water source.

**Section 3 - Summary of Source Water Assessment and Contaminant Assessment (DDW Step 1):** This section comprises a summary discussion of the Step 1 results for the Olympic Well Field.

**Section 4 - RWQC Data:** This section describes the data and data sources used for the Olympic Well Field RWQC study.

**Section 5 - RWQC Approach:** This section describes the approach adopted for characterization of raw water quality for the Olympic Well Field.







**Section 6 - RWQC Results:** This section presents the results of the RWQC study.

**Section 7 - Conclusions:** This section provides a summary of the results and conclusions of the RWQC study.

**Section 8 - References:** This section lists the documents referenced in this RWQC report.







# 2 Olympic Well Field Water Source

This section presents a description summary of the Olympic Well Field water source. A detailed discussion is provided in the *Olympic Well Field Drinking Water Source Assessment and Contaminant Assessment* report (Step 1 of 97-005 Evaluation, ICF 2020).

The Olympic Well Field extracts water from the underlying local and regional aquifers, which predominantly comprises permeable sands and gravels interbedded with laterally discontinuous lenses of less permeable finer-grained silt and clays. Locally, the upper water-bearing zones in the Olympic Well Field are identified as the B-, C-, and D-Zones. In the regional context, the B-Zone is known as the Lakewood Aquifer and C- and D-Zones are within the Silverado Aquifer. The C-Zone within the Olympic Well Field is laterally continuous, whereas the B- and D-Zones pinch-out in the western portion of the well field. The Sunnyside Formation, a regionally recognized aquifer, underlies the D-Zone. Production wells SM-3 and SM-4 were installed in 1969 and 1981, respectively, before contamination was discovered in the Olympic Well Field. Each was screened in several water-bearing units, including the B- and C-Zones, the zones most impacted by contamination. An idealized geological cross-section with local to regional aquifer correlation is shown in Figure 2-1.

Static (non-pumping) water levels in Olympic Well Field production wells are approximately 100 feet below ground surface (ft bgs); dynamic water levels in pumping wells generally vary between 250 and 300 ft bgs, depending on pumping rates. Under natural conditions (non-pumping) groundwater flows west toward the Pacific Ocean. The depth to water in monitoring wells in the B- and C-Zones generally ranges between 100 and 120 ft bgs and 150 to 170 ft bgs, respectively. However, water levels can extend beyond those ranges depending on several factors, including pumping rates in production wells and seasonal variations in recharge. Locally, groundwater hydraulic gradients vary in magnitude and direction depending on various groundwater stresses, such as groundwater pumping and limited recharge during drought conditions.

A calibrated groundwater flow model developed by ICF for the Olympic Subbasin (ICF 2017) was used to simulate the future capture zones generated by pumping of production wells SM-4, SM-8 and SM-9 which are planned to supply the treatment system. Capture zones were generated for each aquifer zone intercepted by the production wells, i.e., the B-Zone, C-Zone, D-Zone and upper Sunnyside Aquifer. Aquifer zones are represented by different layers within the groundwater model. The pumping rates specified in the model for each production well were based on planned future pumping rates; these are summarized in Table 2-1 below.

For the purposes of this report, the Olympic Well Field Study Area was delineated by determining an aggregate capture zone comprising the modeled 2-, 5-, and 10-year capture zones for the B-, C-, D-Zones and Sunnyside Aquifer as a means of predicting the lateral extent the groundwater system will be influenced by Olympic Well Field pumping activities. The Olympic Well Field Study Area is shown in Figure 2-2. In addition, modeled particle traces and 2-, 5-, and 10-year capture zones for each aquifer zone are presented in the figures provided in Appendix A.

The forecast (forward-looking) flow model simulation was performed assuming steady-state pumping conditions to assess long-term flow patterns. Simulated particles were then placed around each production well in the model and reverse particle tracking was performed to trace the lateral extent of capture for each well in the applicable aquifer zone where the wells are screened.



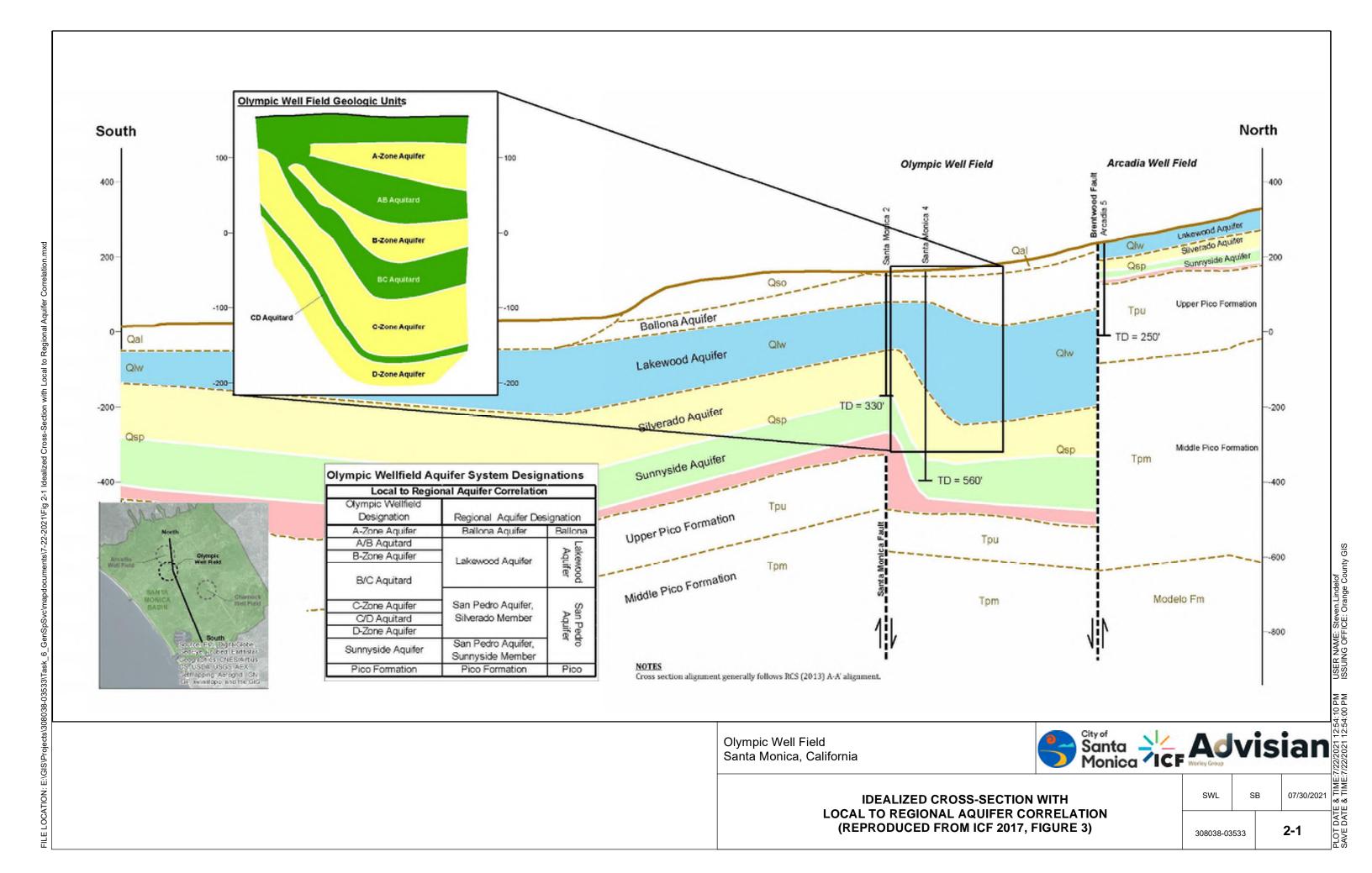


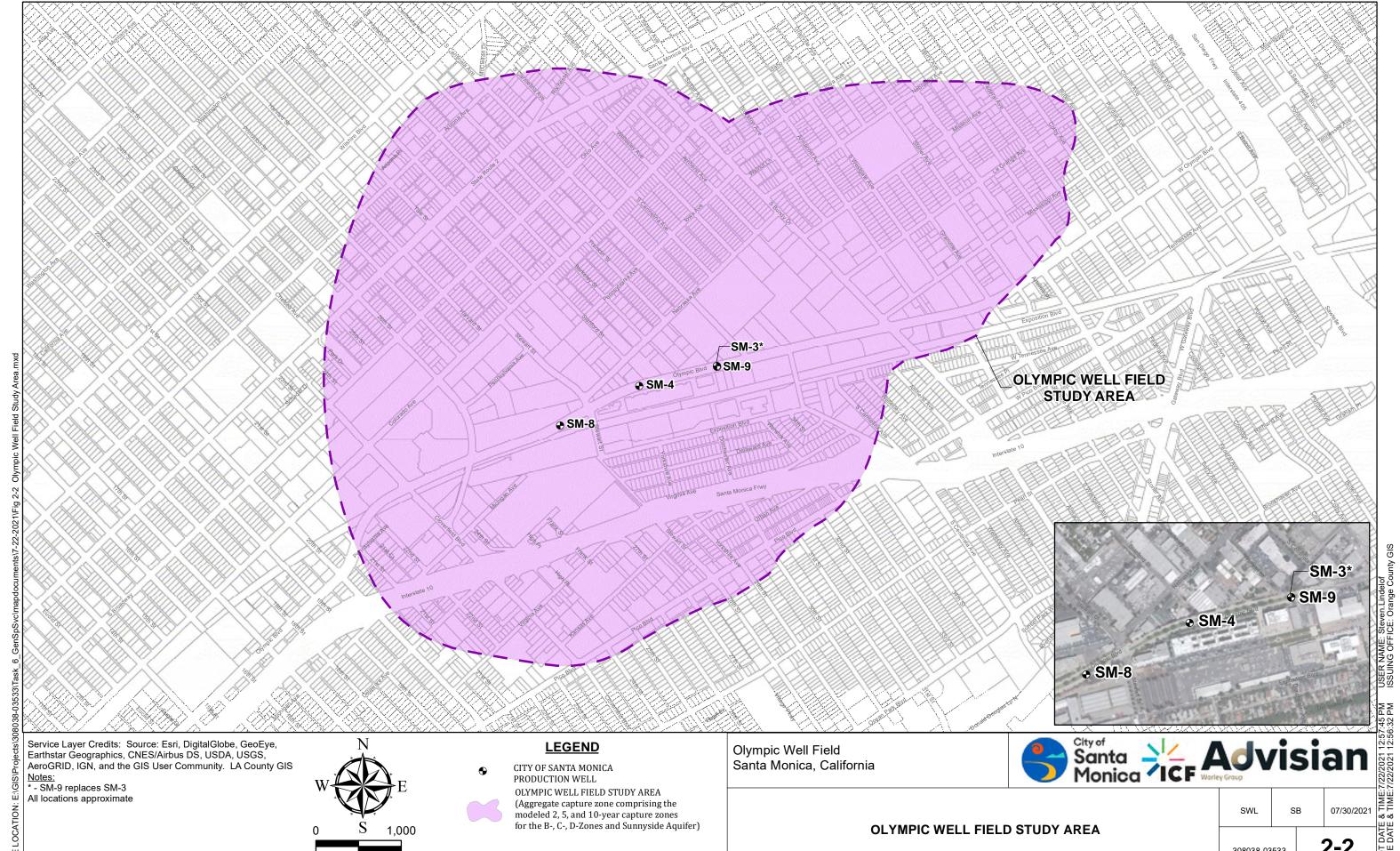


Table 2-1 Modeled Production Well Pumping Rates and Aquifers/Zones Intercepted by Olympic Well Field Production Wells

Production Well	Aquifers/Zones Intercepted	Modeled Future Pumping Rate (gpm)	
	B-Zone		
CN4.4	C-Zone	700	
SM-4	D-Zone	700	
	Sunnyside Aquifer		
	C-Zone		
SM-8	D-Zone	700	
	Sunnyside Aquifer		
	B-Zone		
	C-Zone		
SM-9	D-Zone	600	
	Sunnyside Aquifer		

**Notes:**  $gpm = gallons \ per \ minute.$ 





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# 3 Summary of Source Water Assessment and Contaminant Assessment Results

As stated in the DDW Process Memo 97-005-R2020, the steps (or elements) of the extremely impaired source evaluation process are designed to be sequential in nature. Step 1 of the 97-005 Evaluation, i.e., the SA/CA, determined the extent to which groundwater within the Olympic Well Field is vulnerable to contaminating activities. The RWQC study (Step 2 of 97-005 Evaluation) documented herein, follows on from Step 1 and includes an evaluation of contaminants and contamination sources identified in the SA/CA. A summary is provided below.

The SA/CA included an assessment of known and potential contaminant sources to identify the origin of detected contaminants in groundwater, and potential contamination sources currently or historically present in the source water capture zones.

The Olympic Well Field is located within a corridor formerly occupied by several industrial operations. Some of these operations resulted in release of volatile organic compounds (VOCs) that have impacted groundwater production wells in the City's Olympic Well Field.

Sources of contamination to groundwater have been identified at the CSHV Pen Factory Property (former Paper Mate/Gillette facility) located at 1681 26th Street and the Former Boeing/Douglas Aircraft property, located at 2902 Exposition Boulevard. The locations of these sites are shown in (Figure 3-1). The City continues to monitor the known sites and evaluates other sites that may also contribute to groundwater contamination in this area. A brief summary of operations and chemical releases at these two sites is provided below, and a detailed discussion is provided in the SA/CA report.

# 3.1 CSHV Pen Factory Property (former Paper Mate/Gillette)

Gillette manufactured ballpoint pens under the Paper Mate brand at 1620 26<sup>th</sup> Street, City of Santa Monica, CA from 1957 until 2006 when operations ceased at this facility. The location of the former site is shown in Figure 3-1. Operations that used potentially hazardous materials, or generated materials designated as hazardous wastes, included degreasing, metal plating, molding, nickel plating, ink manufacturing, product assembly, and tooling and maintenance. Over the facility's history, virgin and spent solvents containing PCE and TCE and other wastes were stored in drums at various locations at the facility and later in up to 35 underground storage tanks (USTs) and 45 aboveground storage tanks (ASTs). USTs and ASTs were generally located in two separate locations, along the northern property line and along the southern property line adjacent to Olympic Boulevard. During investigations and closure activities, leaking USTs and associated underground piping were discovered, as well as the presence of contaminated soils in multiple locations at the former facility (AMEC, 2008). As of the late 2000s, all known USTs and ASTs had been removed and numerous investigations and extensive source area remediation had been performed.

Between 2010 and 2015 soil remediation and shallow groundwater remediation occurred at the site to remove PCE and TCE. The Los Angeles Regional Water Quality Control Board (LARWQCB) subsequently concluded that remedial activities had cleaned-up or abated contamination to assure protection of groundwater beneath the site and vicinity for its beneficial uses (LA-RWQCB, 2016). This site is now used for commercial office space and other non-industrial purposes.



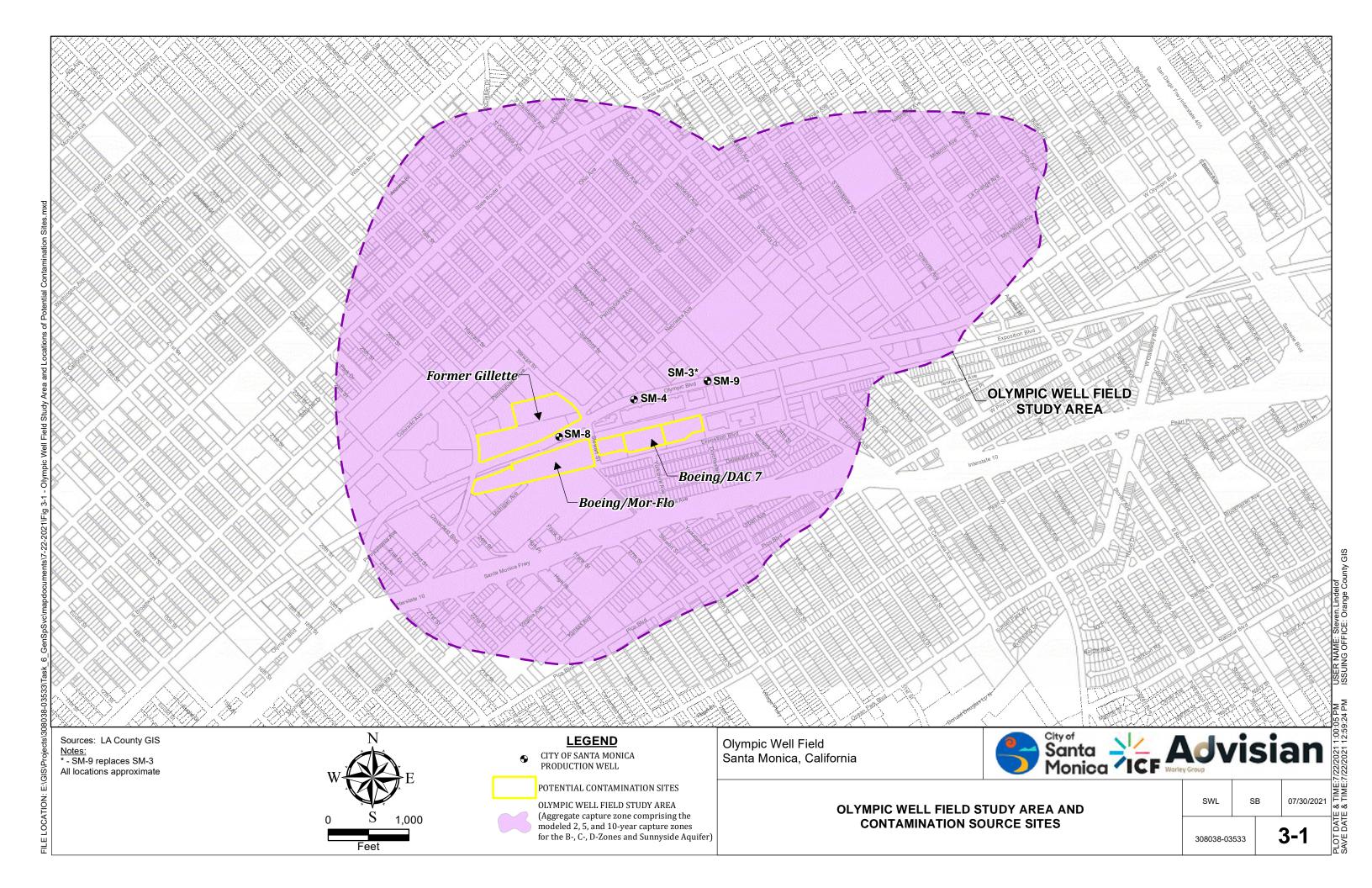




## 3.2 Former Boeing/Douglas Aircraft Property

Douglas Aircraft (Douglas) operated an aerospace components company (Plant A7) located at 2902 Exposition Boulevard and used the eastern and central portions of what is now known as the Bergamot Arts Center for warehousing. The location of the former site is shown in Figure 3-1. Douglas later merged with McDonnell Aircraft and became McDonnell Douglas. In 1997, the Boeing Company acquired the assets of McDonnell Douglas. Aircraft parts manufacturing commenced at the Plant A7 site in the mid-1940s and ceased in the early-1970s. Douglas utilized industrial solvents in their manufacturing process. General Telephone & Electronics Corporation (GTE) acquired the property from Boeing in the early-1970s and operated a fleet refueling and maintenance facility until they sold the property to Verizon in the late 1970s. VOCs, principally PCE and TCE have been detected in the soil, vapor, and shallow groundwater (A-Zone) beneath the Boeing-Exposition Boulevard site. VOC releases to the A-Zone at the Boeing site have been characterized by over 150 soil borings and groundwater monitoring wells.

Verizon utilized the former Plant A7 property for warehousing, vehicle maintenance, and fleet washing beginning in the early 1970s. Releases of fuel hydrocarbons occurred to the soil and shallow groundwater during Verizon's occupancy of the Plant A7 site. No fuel related compounds have been detected in the City's Olympic Well Field production wells. Remedial activities were conducted to remediate the fuel hydrocarbon releases. Notably, Boeing has retained obligations to address shallow groundwater VOC contamination at this site, which includes monitored natural attenuation and deeper groundwater assessment. Much of the former Plant A7 site is now occupied by a Los Angeles County Metropolitan Transportation Authority (Metro) maintenance facility. The rest of the former Plant A7 site is now office space.









# 4 Raw Water Quality Characterization Data

This section provides information pertaining to the analysis period and data used for the data analyses conducted for the RWQC study. Data types include:

- Groundwater monitoring well water quality data;
- Production well data which includes:
  - water quality data;
  - groundwater production data; and,
- Precipitation data.

For the purposes of the RWQC study, an analysis period of January 1, 2012 to June 30, 2020 was selected. The rationale for selecting 2012 as the start of the analysis period is because in-situ soil remediation at one of the primary identified responsible parties<sup>1</sup> in the Olympic Well Field concluded in February 2012. It was considered suitable to use data after completion of remediation as these activities are likely to have influenced constituent concentrations and therefore concentrations recorded prior to completion of remediation are unlikely to be representative of current or future groundwater conditions. The analysis period also includes a range of wet and dry years to evaluate long-term variations in COPC concentrations; this is discussed later in Section 6.5.

The RWQC data sources are described in the following subsections.

# 4.1 Olympic Well Field Monitoring Well Water Quality Data

As described in Section 2, reverse particle tracking was performed to delineate the lateral extent of the projected capture for each production well in the applicable aquifer zone(s) where each well is screened. Once capture zones were generated, groundwater monitoring wells located within these capture zones were identified. The capture zones for each production well and each applicable aquifer zone, as well as the locations of the identified groundwater monitoring wells are shown in the figures provided in Appendix A. A total of 22 groundwater monitoring wells were selected and the locations of these wells are shown in Figure 4-1. Table 4-2 presents the list of monitoring wells, the aquifer zone each monitoring well intercepts, the production well capture zone assigned to each monitoring well, and pertinent well construction information.

Groundwater quality data for an extensive list of constituents (>500) were queried from the Olympic Well Field Environmental Quality Information System (EQuIS) database for the 22 selected groundwater monitoring wells. The database stores water quality data generated from the implementation of the following groundwater monitoring programs at the Olympic Well Field:

• The Olympic Well Field Monitoring Program which comprises quarterly groundwater monitoring activities. The program commenced in 2002 and is still ongoing. The groundwater monitoring results are documented and reported to the LARWQCB. Samples collected as part of this monitoring program

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<sup>&</sup>lt;sup>1</sup> Former Gillette/Paper Mate Site located at 1620 26<sup>th</sup> Street, City of Santa Monica, CA. Refer to Section 3 for more information.







are analyzed using United States Environmental Protection Agency (EPA) Method 8760. However, when using this method, the laboratory method reporting limits (MRLs) for 1,4-dioxane and 1,2,3-TCP are above the NL (1  $\mu$ g/L) and MCL (0.005  $\mu$ g/L), respectively. Therefore, additional low-level analyses are conducted using EPA Method 522 for 1,4-dioxane (MRL = 0.5  $\mu$ g/L) and EPA-approved method CA SRL 524M for 1,2,3-TCP (MRL = 0.005  $\mu$ g/L). Quarterly groundwater monitoring reports are available on GeoTracker (Global ID: T0603799303;

https://geotracker.waterboards.ca.gov/profile report?global id=T0603799303).

- The 97-005 Raw Water Characterization Sampling and Analysis Plan (Advisian 2020) which was executed in 2020 for the purpose of generating data of sufficient quality and quantity to support full characterization of the raw water quality in the Olympic Well Field, meeting DDW requirements. Groundwater samples were analyzed for >500 constituents, including pesticides, pharmaceuticals, radionuclides, tentatively identified compounds (TICs) and other various organic and inorganic compounds. Samples were analyzed using EPA-approved drinking water methods. MRLs are equal to or less than applicable MCLs or NLs.
- Implementation of the Monitoring and Reporting Program (MRP) No. CI-10539 for Santa Monica Exploratory Well SM-10i (LARWQCB 2020). The LARWQCB requires that the City conduct quarterly and annual monitoring of monitoring wells in the receiving aquifer surrounding exploratory well SM-10i which was used to perform short-term injection testing within the Olympic Well Field (documented in MRP). Samples were analyzed using EPA-approved drinking water methods. MRLs are equal to or less than applicable MCLs or NLs. Technical and monitoring reports associated with the MRP is available on GeoTracker (WDR100046020;

https://geotracker.waterboards.ca.gov/profile report.asp?global id=WDR100046020).

The data generated from the implementation of these monitoring programs and exported from the Olympic Well Field EQuIS database, were used to screen for contaminants of potential concern (COPCs) (Section 6.1) and estimating future treatment plant influent concentrations (Section 6.1.3).

# 4.2 Olympic Well Field Production Well Data

The capture zones for each production well and each applicable aquifer zone, as well as the locations of the identified groundwater monitoring wells for each aquifer zone are shown in Figures A-1 to A-4, provided in Appendix A.

The City conducted sampling of production well SM-4, and newly installed, but yet to be equipped and permitted production wells SM-8 and SM-9 (SM-3 replacement well) concurrent with the implementation of the 97-005 Raw Water Characterization Sampling and Analysis Plan (Advisian 2020). The purpose of this sampling was to generate production well data of sufficient quality and quantity to support full characterization of the raw water quality in the Olympic Well Field. Groundwater samples were analyzed for >500 constituents, including pesticides, pharmaceuticals, radionuclides, TICs and other various organic and inorganic compounds. This data was used to identify COPCs (Section 6.1) and for comparisons to estimated future treatment plant influent concentrations derived from monitoring well data (Section 6.1.3).

The City conducts compliance sampling of the Olympic Well Field production wells (SM-3 and SM-4) in accordance with its Domestic Water Supply Permit (DDW 1910146). Samples are collected by the City on a monthly basis; therefore, a significant water quality data set has been established. Water quality data generated from compliance sampling are transmitted in electronic format to DDW. In addition to







collecting water samples, the City also measures and records water level and production data for each production well. Temporal data generated from monthly compliance monitoring were used to evaluate water quality concentration trends and variability (Section 6.4).

The locations of the production wells are shown in Figure 4-1 and pertinent well information is presented in Table 4-3.

## 4.3 Precipitation Data

Annual precipitation data was obtained from Santa Monica Municipal Airport (Station Number: USW00093197) for the January 1, 2012 through June 30, 2020 period and is presented in Table 4-1. This data was used determine whether changes in water quality could be correlated to periods considered to be "wet" (i.e., periods of high precipitation) versus "dry" (i.e., periods of low precipitation) (Section 6.5).

Table 4-1 Annual Precipitation at Santa Monica Municipal Airport

Calendar Year	Water Year (Start Date)	Total Precipitation (inches)
2019-2020	1-Oct-19	13.68
2018-2019	1-Oct-18	19.81
2017-2018	1-Oct-17	5.93
2016-2017	1-Oct-16	18.09
2015-2016	1-Oct-15	6.43
2014-2015	1-Oct-14	11.66
2013-2014	1-Oct-13	4.76
2012-2013	1-Oct-12	7.24

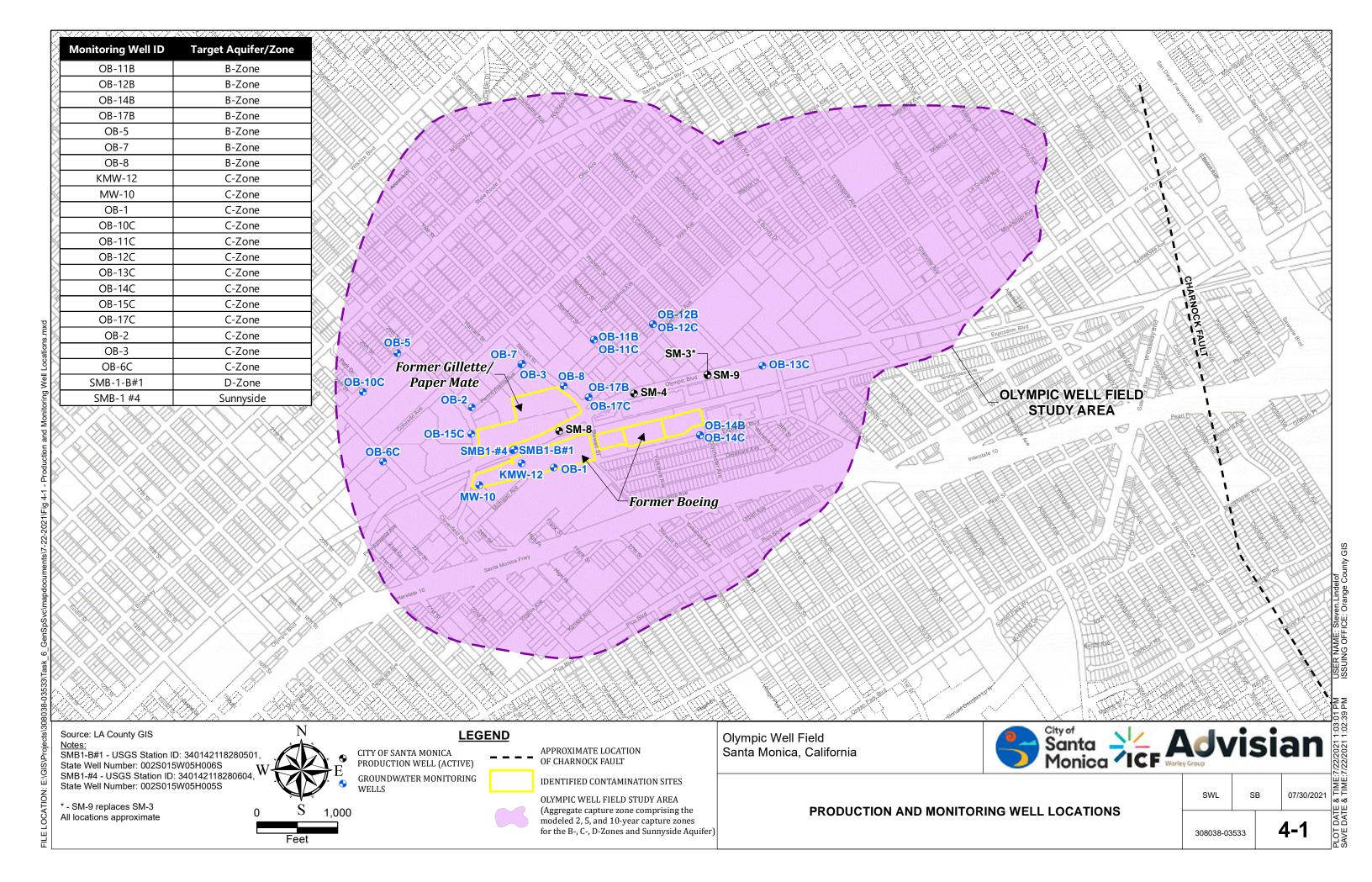






Table 4-2 Olympic Well Field Monitoring Wells Information

W. II ID	W.II.O	A	Production Well	Footbal	News	Ground Surface	Top of Casing	Screen Interval (ft bgs)		Screen Interval (ft AMSL)	
Well ID	Well Owner	Aquifer Zone	Capture Zone	Easting	Northing	Elevation (ft AMSL)	Elevation (ft AMSL)	Тор	Bottom	Тор	Bottom
KMW-12	City	C-Zone	SM-8	6419868.9	1832919.6	156.2	155.5	140.0	165.0	16.2	-8.8
MW-10	City	C-Zone	SM-8	6419343.6	1832650.7	153.4	155.6	123.5	143.5	29.9	9.9
OB-1	City	C-Zone	SM-8	6420268.4	1832870.3	156.3	156.0	139.8	159.8	16.5	-3.5
OB-10C	City	C-Zone	SM-8	6417901.4	1833803.9	152.4	151.7	168.3	198.3	-15.9	-45.9
OB-11B	City	B-Zone	SM-4	6420765.9	1834448.6	156.7	156.2	230.2	250.2	-73.5	-93.5
OB-11C	City	C-Zone	SM-4	6420765.9	1834448.6	156.7	156.0	346.3	376.3	-189.6	-219.6
OB-12B	City	B-Zone	SM-4	6421494.2	1834642.3	156.8	156.3	220.3	240.3	-63.5	-83.5
OB-12C	City	C-Zone	SM-4	6421494.2	1834642.3	156.8	156.3	336.5	366.5	-179.7	-209.7
OB-13C	City	C-Zone	SM-9	6422849.8	1834130.3	154.2	153.9	305.6	335.6	-151.4	-181.4
OB-14B	City	B-Zone	SM-9	6422077.2	1833265.5	150.8	150.5	102.0	132.0	48.8	18.8
OB-14C	City	C-Zone	SM-9	6422077.2	1833265.5	150.8	150.4	186.4	216.4	-35.6	-65.6
OB-15C	City	C-Zone	SM-8	6419246.2	1833281.4	156.1	156.1	185.0	215.0	-28.9	-58.9
OB-17B	City	B-Zone	SM-4	6420698.9	1833740.4	152.6	152.6	194.9	204.9	-42.3	-52.3
OB-17C	City	C-Zone	SM-4	6420698.9	1833740.4	152.5	152.5	295.6	325.6	-143.1	-173.1
OB-2	City	C-Zone	SM-8	6419251.0	1833613.0	155.2	154.9	246.5	276.5	-91.4	-121.4
OB-3	City	C-Zone	SM-8	6419874.2	1834156.2	157.1	157.5	339.3	369.3	-182.2	-212.2
OB-5	City	B-Zone	SM-4	6418328.2	1834284.6	156.0	155.7	145.9	175.9	10.1	-19.9
OB-6D	City	C-Zone	SM-8	6418153.1	1832939.5	149.2	148.8	145.9	175.9	3.3	-26.7
OB-7	City	B-Zone	SM-4	6419865.3	1834147.6	157.5	157.1	215.0	246.0	-57.6	-88.6
OB-8	City	B-Zone	SM-4	6420392.0	1833878.0	153.8	153.4	205.0	235.0	-51.2	-81.2
SMB1-B#1 <sup>(1)</sup>	USGS	D-Zone	SM-4, SM-8, SM-9 <sup>(3)</sup>	6419779.1	1833088.0	159.1	159.3	235.0	255.0	-75.9	-95.9
SMB1-#4 <sup>(2)</sup>	USGS	Sunnyside	SM-4, SM-8, SM-9 <sup>(4)</sup>	6419759.9	1833082.6	159.3	159.1	350.0	370.0	-190.7	-210.7

**Notes:** ft = feet; AMSL = Above Mean Sea Level; bgs = below ground surface; City = City of Santa Monica; USGS = United States Geological Survey

<sup>(1)</sup> SMB1-B#1 = USGS Station ID: 340142118280501; State Well Number: 002S015W05H006S

<sup>(2)</sup> SMB1-#4 = USGS Station ID: 340142118280604; State Well Number: 002S015W05H005S

<sup>(3)</sup> SMB-1-B#1 used for all production well capture zones in D-Zone

<sup>(4)</sup> SMB-1 #4 used for all production well capture zones in Sunnyside Aquifer







Table 4-3 Olympic Well Field Production Wells Information

Well Name	Easting	Northing	Year Drilled	Ground Surface Elevation (ft AMSL)	Pump Intake Depth (ft bgs)	Elevation of Suction Intake (ft AMSL)	Perforation Interval Depths (ft bgs)	Perforation Interval Elevations (ft AMSL)
SM-3	64222173.0	1834009.0	1969	155.0	280.0	-125.0	210 to 270 300 to 380 410 to 430 490 to 530	-55 to -115 -145 to -225 -255 to -275 -335 to -375
SM-4	6421262.9	1833785.3	1981	150.7	240.0	-89.3	200 to 410 470 to 540	-49 to -259 -319 to -389
SM-8	6420334.0	1833323.0	2017	156.0			210 to 265 295 to 325 335 to 345 360 to 460	-54 to -109 -139 to -169 -179 to -189 -204 to -304
SM-9	6422172.63	1834014.79	2019	155.0			240 to 265 300 to 380 390 to 430 490 to 535 655 to 750 760 to 790	-85 to -110 -145 to -225 -235 to -275 -335 to -380 -500 to -595 -605 to -635

**Notes:** ft = feet; AMSL = Above Mean Sea Level; bgs = below ground surface, -- pump not installed at time of writing this report



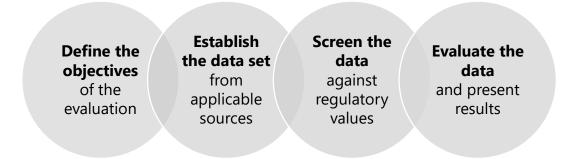




# 5 Raw Water Quality Characterization Approach

A series of steps were taken in order to evaluate data to support this RWQC; these steps are summarized in Figure 5-1. This section outlines the approach and methodologies adopted for each step of the evaluation. The results of the evaluation are presented later in Section 6.

Figure 5-1 Key Steps of the Evaluation



## 5.1 Define the Objectives

Per the DDW Process Memo 97-005-R2020 (DDW 2020), the primary objective of the RWQC is to characterize the influent water quality that will enter the treatment system, so that an appropriate level of monitoring and treatment can be designed to produce safe drinking water. This includes an evaluation of potential future water quality at the production wells and plant influent for the contaminants found in Step 1 of the 97-005 Evaluation, i.e., the SA/CA (ICF 2020; refer to Section 3).

#### 5.2 Establish the Data Set

From the sources described in Section 4, water quality sample records for production and monitoring wells were combined into a single data set. The water quality data set was then queried to identify constituents analyzed in groundwater well samples using specific criteria as follows:

- Well type (production or monitoring well);
- Well screen depth/elevation interval and associated aquifer zones intercepted by the well screen;
- Well location within modeled captures zones for each aquifer zone; and,
- Date of samples (samples collected between January 1, 2012 and June 30, 2020<sup>2</sup>).

The resultant data set consisted of 48,151 records from over 650 samples, featuring over 500 constituent/parameter results derived from four production wells and 22 monitoring wells.

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<sup>&</sup>lt;sup>2</sup> An exception is 1,4-dioxane and 1,2,3-trichloropropane (1,2,3-TCP) data for USGS wells SMB1-B#1 and SMB1-#4; data were not available for these constituents for the time period, so groundwater samples collected and analyzed for these constituents in Q3 2020 were used for the evaluation.







#### 5.3 Screen the Data

To identify COPCs and determine which constituents would drive design for proposed treatment, the groundwater quality data set (Section 5.2) was screened against specific criteria. This section discusses the approach adopted for COPC screening for this RWQC, with the results being presented later in Section 6.

Starting with the entire data set, an initial COPC screening was conducted by applying the following criteria to develop a preliminary list of COPCs for further evaluation:

- 1. Constituents that were detected in the four production wells (SM-3, SM-4, SM-8, or SM-9);
- 2. Constituents that were non-detect in the four production wells, but had three or more detections in one or more monitoring wells;
- 3. Detected constituents with a primary maximum contaminant level (MCL) or notification level (NL); and,
- 4. Constituents meeting screening criteria 1 or 2 and screening criteria 3 with a ratio of detected concentration (maximum observed) to MCL or NL greater than 0.05 (5%).

Following the initial screening, the resultant list of COPCs were further refined to determine which COPCs would drive design for the planned treatment system, using the following secondary criteria:

- 5. Constituents meeting initial screening criteria with a ratio of detected concentration (maximum observed) to MCL or NL greater than 0.5 (50%); and,
- 6. Constituents meeting initial screening criteria and screening criteria 5 which are synthetic organics (these constituents are most likely to drive design for the planned treatment system using ultraviolet/advanced oxidation process [UV/AOP] and granular activated carbon [GAC]).

Results from the initial and secondary screening are provided in Section 6.1.

#### 5.4 Evaluate the Data

The next stage in the RWQC was to evaluate the water quality data. As part of the evaluation, several different data analyses were conducted to meet the requirements of the DDW Process Memo 97-005-R2020 (DDW 2020), including:

- Estimating future treatment plant influent concentrations to characterize the quality of water anticipated to enter the proposed treatment system;
- Trend analysis for historical water quality to understand how water quality trends have changed over time; and,
- Variability/correlation analysis to develop an understanding of how water quality has changed over time under the influence of certain factors such as pumping and seasonal variation in precipitation.

The approach adopted for each of the data analyses are described in the subsections which follow.

## 5.4.1 Approach to Estimating Future Treatment Plant Influent Concentrations

The approach to estimating future treatment plant influent concentrations utilizes several methods including statistical analysis to characterize ranges in raw water quality in the data set and, subsequently,







flow-weighting calculations to estimate concentrations in production wells and treatment plant influent. This section discusses the approach adopted for estimating future treatment plant influent concentrations for this RWQC, with the results being presented later in Section 6.

#### 5.4.1.1 Statistical Analysis of Groundwater Quality Data Set

The approach to the COPC screening evaluation (Section 5.3) utilized data from both production and monitoring wells to provide a conservative first stage in characterizing COPCs. However, in order to provide projections of future water quality in production wells and treatment plant influent, monitoring well data were primarily used since these data represent groundwater that is anticipated to arrive at the production wells in the future based on groundwater modeling and capture zone analysis (refer to Section 2).

Using the monitoring well groundwater quality data set, statistical analysis was performed using the EPA's ProUCL software version 5.1 (EPA 2015a, 2015b). ProUCL is a comprehensive statistical software package with statistical methods to evaluate environmental sampling data. For the purposes of the evaluation herein, the mean and 95 percent (%) upper confidence limit of the population mean (UCL95) were calculated using ProUCL software. In the event either of these statistics could not be calculated (for example, due to an insufficient number of detections to compute reliable and meaningful statistics and estimates), the maximum observed value was adopted in subsequent calculations as a conservative approach.

The statistical analysis was conducted for data from monitoring wells within the projected capture zone for each production well supplying the treatment plant (SM-4, SM-8, SM-9) and for each aquifer zone which the production wells are screened in. The results of this analysis are provided in Section 6.1.2.

### 5.4.1.2 Flow-Weighting Calculations for Production Well Concentration Estimates

Following statistical analyses of the monitoring well groundwater quality data set, constituent concentrations based on the mean and UCL95 of the data set were used in flow-weighting calculations to derive production well concentration estimates, applying the following equation:

$$C_w = \frac{c_1 Q_1 + c_2 Q_2 + c_3 Q_3 + c_4 Q_4}{Q_w}$$
 Equation (1)

Where:

C<sub>w</sub> is the estimated flow-weighted concentration in the production well.

 $C_1$  to  $C_4$  is the assumed concentration within each capture zone and within each aquifer zone (adopted from mean or UCL95 statistics, or maximum observed value if either of these statistics were unavailable).

 $Q_w$  is the total flow rate for the production well.

 $Q_1$  to  $Q_4$  is the flow rate to the well in each aquifer zone. This was estimated by calculating the proportion of total flow in the production well coming from each screen interval using the transmissivity across each production well screen within each aquifer zone. Transmissivities were







calculated by multiplying total screen length in each aquifer zone by the hydraulic conductivity for the applicable layer from the calibrated groundwater model.

It should be noted that this approach is conservative as it assumes there are no transport mechanisms that reduce constituent concentrations, such as attenuation, retardation or degradation. Furthermore, no monitoring network is robust enough to describe all subsurface conditions; the approach herein assumes all flow from each aquifer zone into each production well screen has concentrations equal to the adopted value from monitoring well statistics (mean or UCL95 statistics, or maximum observed value if either of these statistics were unavailable). This may result in an over-estimate of production well influent concentrations for some constituents as there are expected to be areas within each capture zone which have low or non-detect concentrations of the constituents evaluated. The results of this analysis are provided in Section 6.1.2.

# 5.4.1.3 Flow-Weighting Calculations for Treatment Plant Influent Concentration Estimates

After concentration estimates for each production well were calculated, plant influent concentrations (comprising the combined flow from the three production wells) were estimated using the same flow-weighting method applied in Equation 1 above, as follows:

$$C_i = \frac{c_{SM4}Q_{SM4} + c_{SM8}Q_{SM8} + c_{SM9}Q_{SM9}}{Q_i}$$
 Equation (2)

Where:

C<sub>i</sub> is the estimated flow-weighted influent concentration for the combined flow of all three production wells (i.e., SM-4, SM-8 and SM-9).

 $C_{SM4}$ ,  $C_{SM8}$  and  $C_{SM9}$  are the estimated concentrations for each production well, derived from Equation 1.

Qi is the combined flow of all three production wells (i.e., SM-4, SM-8 and SM-9).

Q<sub>SM4</sub>, Q<sub>SM8</sub> and Q<sub>SM9</sub> are the flow rates for each production well.

For treatment plant influent estimates, only the production well estimates derived from UCL95 statistics (or maximum observed value if these statistics were unavailable) were used as this is more conservative than the mean. It should be noted that the flow-weighted treatment plant influent concentration estimates may, in some cases, be significantly lower than any single observed value in the groundwater monitoring well data set. The results of this analysis are provided in Section 6.1.2.

## 5.4.2 Trend Analysis

Trend analysis for historical water quality data provides important information that supports the evaluation of future COPC concentrations in the production wells to help understand how water quality trends have changed over time. For example, increasing trends indicate that constituent concentration levels have not stabilized, and concentrations may continue to increase to levels greater than historically and currently observed. This information supports proper treatment design as it provides insight to potential future







influent concentrations. This section discusses the approach adopted for trend analysis for this RWQC, with the results being presented later in Section 6.

Available concentration data for COPCs between January 1, 2012 and June 30, 2020 were used for trend analysis. Trend analysis was undertaken for water quality data from production wells SM-3 and SM-4 to assess concentration trends and evaluate temporal changes in COPCs. These production wells were chosen because extensive temporal data exist for SM-3 and SM-4 but limited temporal data is available for the recently drilled production wells SM-8 and SM-9 (refer to Section 2). As noted in Section 4.2, SM-9 was drilled as a replacement well for SM-3. SM-3 is screened across four depth intervals from 210 to 530 ft bgs, whereas SM-9, located within approximately 50 ft of SM-3, is screened over six depth intervals, from 240 to 790 ft bgs. The upper four screen intervals of SM-9 align very closely with the screen intervals of SM-3, and therefore it is reasonable to use historic data from SM-3 as a proxy for historical water quality data for SM-9. In addition, the two lower screen intervals of SM-9 (between 655 and 790 ft bgs) draw water from deeper aquifer units that are likely to be uncontaminated, or less contaminated, than the shallower four screen intervals, and therefore is anticipated to dilute contaminant concentrations in SM-9 compared to SM-3. Thus, using SM-3 as a proxy for SM-9 data is likely to be conservative.

#### 5.4.2.1 Methodology

Trend analysis on chemical data was performed using the non-parametric Mann-Kendall Trend Test (Kendall 1975; Mann 1945), and a modified version of Sen's non-parametric slope estimator (Sen 1968). The process involves first arranging the data in a chronological order. The sign of the differences between all forward combinations of data pairs are then used to compute the Mann-Kendall statistic (S). Next, the Mann-Kendall S and the number of data points (n) are used for the normal approximation test. The results yield a probability value from 0 to 1.

The Sen's slope was used to provide an approximate magnitude of trends. Sen's slope is calculated by evaluating the slope between all forward combinations of data pairs. The slopes are then ranked, and the median slope value is selected. Unlike least squares regression, the Sen's slope is not greatly affected by gross outliers, which are occasionally found in groundwater chemistry data. Slopes are presented in absolute terms (µg/L per yr.) and normalized relative to the median concentration (% change per yr.).

The trend analysis was conducted for two time periods to evaluate the most recent water quality trends:

- January 2012 to June 2020 data series; and,
- January 2017 to June 2020 data series.

Trend analysis of the 2012 to 2020 data series provides a longer-term assessment of temporal changes in water quality, the shorter duration 2017 to 2020 time period allows for a more current evaluation of COPC trends.

Trend analysis data were screened according to the following three criteria to indicate potentially significant temporal trends:

1. The probability of a temporal trend - an inferred confidence level greater than 0.90 was used for the first level of screening. This corresponds with the general convention that a 90% confidence interval is often used in statistical reporting.







- 2. Normalized value of Sen's slope indicator (% change per year) a normalized slope threshold of >5% per year was used to screen out relatively small trends superimposed on high parameter concentrations.
- 3. Absolute value of Sen's slope indicator (µg/L/yr) calculated and used to screen out high probability trends with very low slopes, based on slope thresholds indicated in Table 5-1. The thresholds were selected based on detection limits and regulatory thresholds. Note that absolute slope criteria are only present for parameters that have sufficient data for analysis (at least eight data points exist and only where more than 50% of the data are above the detection limits).

Table 5-1 Absolute Slope Thresholds of Constituents of Potential Concern, Sen's Slope Method

СОРС	Absolute Slope Criteria (μg/L/year)				
1,1-DCE	± 0.5				
1,4-Dioxane	± 0.1				
cis-1,2-Dichloroethene	± 0.5				
Nitrate and Nitrite as Nitrogen (as N)	± 200				
PCE	± 0.5				
TCE	± 0.5				
Total Trihalomethanes	± 5				

#### 5.4.2.2 Limitations of Statistical Trend Approach

The Mann-Kendall and Sen's Slope trend calculations assume that:

- Data is monotonic (i.e., entirely increasing or entirely decreasing); and,
- In the absence of a trend, the individual data points within any given data set are independent of each other (i.e., no serial correlations where the level of a variable affects its future level).

The Mann-Kendall and Sen's Slope trend calculations were applied to the SM-3 and SM-4 production well data sets. Additionally, a visual assessment of trends was conducted to identify cases where data does not meet the assumptions of the Mann Kendall and/or Sen's slope calculations. A single data set could contain multiple trends, trend reversals, stabilization, and/or newly developing trends due to several factors including:

- Complexities in contaminant distribution(s);
- Arrival of a new contaminant plume at a well;
- Contaminant plume core(s) passing by a well; and/or,
- Concentration variations due to activation and deactivation of groundwater production at a well or nearby well(s).







The purpose of the visual trend assessment review is to provide a qualitative assessment of the current concentration trends in the context of these complexities which is incorporated into the discussion of results. The results of trend analyses are provided in Section 6.4.

#### 5.4.3 Raw Water Quality Variability

Evaluating the variability of water quality data helps develop an understanding of how water quality has changed over time under the influence of certain factors such as pumping of production wells. In accordance with the DDW Process Memo 97-005 -R2020 (DDW 2020), an assessment of COPC concentration variability with pumping rate and time (seasonal and long-term) observed in the production wells was conducted to inform the understanding of future potential variability in concentrations that may arrive at the planned treatment plant. This section discusses the approaches adopted to evaluate raw water quality variability for this RWQC, with the results being presented later in Section 6.

#### 5.4.3.1 COPC Concentration Variations with Pumping Rate

The analysis focuses on the observed relationships between production (pumping) at a single production well and the concentration of each COPC using water quality data from production wells SM-3 and SM-4. Only COPCs that contained sufficient temporal resolution during the analysis period (January 1, 2012 to June 30, 2020) are assessed. Trend charts are developed and used to assess potential concentration changes with respect to pumping rates. The data are visually assessed to determine correlations between COPC concentrations and pumping, with a positive correlation being defined as instances where the recorded concentration of a parameter increases (or decreases) as pumping rate increases (or decreases). On the other hand, a negative correlation is defined as instances where the recorded concentration of a parameter increases (or decreases) as the pumping rate decreases (or increases). The results of this analysis are provided in in Section 6.

#### 5.4.3.2 COPC Concentration Variations with Time (Seasonal and Wet / Dry Period)

The analysis focuses on the observed relationships and patterns between different seasons and concentration of each COPC using water quality data from production wells SM-3 and SM-4. This analysis aims to determine if seasonal impacts or seasonal signals are distinguishable in the data sets.

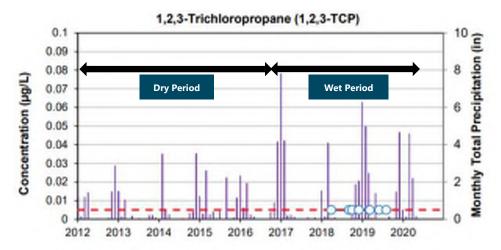
Variability of the climate over time was also assessed with respect to changes in COPC concentrations. The objective of this evaluation was to determine whether changes in concentrations of COPCs could correlated to periods considered to be "wet" (i.e., periods of high precipitation) versus "dry" (i.e., periods of low precipitation). With reference to Figure 5-2 below, the dry period is represented by years 2012-2016, and the wet period 2017-2020. The results of this analysis are provided in in Section 6.







Figure 5-2 Example Chart: Production Well SM-3 – 1,2,3-TCP Concentration versus Precipitation (City of Santa Monica Airport)









# 6 Raw Water Quality Evaluation Results

This section presents the results of each step of the raw water quality evaluation using the approach and methodologies outlined previously in Section 5.

#### 6.1 Screening

#### 6.1.1 COPC Identification

Based on the initial screening (using criteria 1 to 4 listed in Section 5.3), 42 COPCs were identified. From the secondary screening (using criteria 5 and 6 listed in Section 5.3), 15 synthetic organic COPCs were identified, with nine from production well data and six from monitoring well data. These are summarized as follows:

- From production well data:
  - 1,1-dichloroethane (1,1-DCA)
  - 1,1-dichloroethene (1,1-DCE)
  - 1,2,3-trichloropropane (1,2,3-TCP)
  - 1,4-dioxane
  - carbon tetrachloride (CTC)
  - cis-1,2-dichloroethene (cis-1,2-DCE)
  - perfluorooctanoic acid (PFOA)
  - tetrachloroethene (PCE)
  - trichloroethene (TCE)

- Additional from monitoring well data:
  - 1,1,2-trichloroethane (1,1,2-TCA)
  - 1,2-dichloroethane (1,1-DCA)
  - benzene
  - methyl tert-butyl ether (MTBE)
  - trans-1,2-dichloroethene (trans-1,2-DCE)
  - vinyl chloride

Screening results for the 42 COPCs identified in the initial screening and the 15 COPCs identified in the secondary screening are provided in Table 6-1. Screening results for all constituents in the water quality data set are provided in Appendix B.





Table 6-1 Summary of COPC Screening

						Pro	duction Wells			Mor	nitoring Wells		Me	eets CO	PC Scr	reening	Criteri	ia*
CASRN	Parameter	Units	MCL	NL	Number of	Number of	Maximum	Ratio of Maximum	Number of	Number of	Maximum	Ratio of Maximum		Init			Secon	
<b>37 13 11 1</b>	. arameter				Observations	Detections	Observed Value	to MCL or NL	Observations	Detections	Observed Value	to MCL or NL	1	2	3	4	5	6
79-00-5	1,1,2-Trichloroethane	μg/L	5	-	34	0	<0.5	-	475	8	11	2.20		✓	✓	1	1	✓
75-34-3	1,1-Dichloroethane	μg/L	5	-	201	2	0.16	0.03	475	67	7.1	1.42	✓	✓	✓	1	✓	✓
75-35-4	1,1-Dichloroethene	μg/L	6	-	201	117	2.9	0.48	475	138	32	5.33	✓	✓	✓	✓	✓	✓
96-18-4	1,2,3-Trichloropropane	μg/L	0.005	-	38	(17)	0.01	2.00	609	29	0.19	38.00	✓	✓	✓	✓	✓	✓
107-06-2	1,2-Dichloroethane	μg/L	0.5	-	201	0	<0.5	-	475	21	6.6	13.20		✓	✓	✓	✓	✓
123-91-1	1,4-Dioxane	μg/L	-	1	68	64	26	26.00	413	367	440	440.00	✓	✓	✓	✓	✓	✓
7429-90-5	Aluminum	μg/L	1000	-	12	7	2400	2.40	38	38	2800	2.80	✓	✓	✓	✓	✓	
7440-36-0	Antimony	μg/L	6	-	12	6	0.2	0.03	38	37	0.39	0.07	✓	✓	✓	✓		1
7440-38-2	Arsenic	μg/L	10	-	12	11	2	0.20	41	39	38	3.80	✓	✓	✓	✓	✓	1
7440-39-3	Barium	μg/L	1000	-	12	12	76	0.08	38	38	100	0.10	✓	✓	✓	✓		1
7440-39-3	Barium	μg/L	1000	-	12	12	76	0.08	38	38	100	0.10	✓	✓	✓	✓		1
71-43-2	Benzene	μg/L	1	-	201	0	<0.5	-	475	24	19	19.00		✓	✓	✓	✓	✓
117-81-7	bis(2-ethylhexyl)phthalate	μg/L	4	-	18	0	<30	-	38	5	0.91	0.23		✓	✓	✓		
7440-42-8	Boron	μg/L	-	1000	7	7	200	0.20	43	41	590	0.59	✓	✓	✓	<b>✓</b>	✓	1
15541-45-4	Bromate	μg/L	10	-	3	0	<25	-	19	6	160	16.00		✓	✓	✓	✓	
56-23-5	Carbon Tetrachloride	μg/L	0.5	-	201	2	0.15	0.30	475	8	1	2.00	✓	✓	<b>✓</b>	✓	✓	✓
7440-47-3	Chromium	μg/L	50	-	13	8	7.2	0.14	38	31	290	5.80	<b>\</b>	✓	<b>\</b>	✓	✓	
156-59-2	cis-1,2-Dichloroethene	μg/L	6	-	201	110	3.4	0.57	475	93	300	50.00	✓	✓	✓	✓	✓	✓
16984-48-8	Fluoride	μg/L	2000	-	9	9	390	0.20	19	19	370	0.19	✓	✓	✓	✓		
12587-46-1	Gross Alpha	pci/L	15	-	7	6	11	0.73	19	19	7.94	0.53	✓	✓	✓	✓	✓	1
118-74-1	Hexachlorobenzene	μg/L	1	-	17	0	<1	-	57	3	0.09	0.09		✓	✓	✓		
7439-92-1	Lead	μg/L	15	-	12	6	4	0.27	32	14	1.2	0.08	✓	✓	✓	✓		
7439-96-5	Manganese	μg/L	-	500	13	13	67	0.13	38	38	340	0.68	✓	✓	✓	✓	✓	
1634-04-4	Methyl tert-butyl ether	μg/L	13	-	200	0	<3	-	475	18	67	5.15		✓	✓	✓	✓	✓
75-09-2	Methylene Chloride	μg/L	5	-	201	0	<0.5	-	475	6	2	0.40		✓	✓	✓		
7440-02-0	Nickel	μg/L	100	-	12	6	4.1	0.04	38	38	200	2.00	✓	✓	✓	✓	✓	
14797-55-8	Nitrate as N	μg/L	10000	-	45	45	7400	0.74	21	17	16000	1.60	✓	✓	✓	✓	✓	
14797-65-0	Nitrite as N	μg/L	1000	-	9	0	<100	-	21	10	160	0.16		✓	✓	✓		
55-18-5	n-Nitrosodiethylamine	μg/L	-	0.01	3	1	0.0015	0.15	19	1	0.0011	0.11	✓		✓	✓		
NN	Nitrate plus nitrite as N	μg/L	10000	-	7	7	5800	0.58	21	17	16000	1.60	✓	✓	✓	<b>_</b>	✓	
14797-73-0	Perchlorate	μg/L	6	-	9	1	1.6	0.27	19	8	6.7	1.12	✓	✓	✓	<b>_</b>	✓	
335-67-1	PFOA	μg/L	-	0.0051	8	6	0.0029	0.57	16	11	0.0033	0.65	✓	✓	✓	✓	✓	✓
1763-23-1	PFOS	μg/L	-	0.0065	8	2	0.0017	0.26	16	1	0.0042	0.65	✓		✓	<b>✓</b>	✓	<b></b>
7782-49-2	Selenium	μg/L	50	-	12	10	4.6	0.09	38	34	9	0.18	✓	✓	✓	✓	$\longrightarrow$	ļ
10098-97-2	Strontium-90	pci/L	8	-	3	3	0.255	0.03	19	19	0.461	0.06	✓	✓	✓	✓	igcup	ļ
127-18-4	Tetrachloroethene	μg/L	5	-	207	205	54	10.80	476	288	470	94.00	✓	✓	✓	✓	✓	✓
THM	Total Trihalomethanes	μg/L	80	-	34	33	4.8	0.06	19	13	18	0.23	✓	✓	✓	✓		
156-60-5	trans-1,2-Dichloroethene	μg/L	10	-	201	0	<0.5	-	475	8	15	1.50		✓	✓	✓	✓	✓
79-01-6	Trichloroethene	μg/L	5	-	205	200	63.2	12.64	475	227	370	74.00	✓	✓	✓	✓	<b>1</b>	✓
7440-61-1	Uranium Rad	pci/L	20	-	7	7	14	0.70	19	19	14	0.70	✓	✓	✓	✓	✓	
7440-62-2	Vanadium	μg/L	-	50	10	10	14	0.28	26	25	11	0.22	✓	✓	✓	✓		
75-01-4	Vinyl Chloride	μg/L	0.5	-	34	0	<0.5	-	475	12	32	64.00		✓	✓	✓	✓	✓

**Notes:** Constituents meeting initial and secondary screening criteria are highlighted **bold** with blue shading; CASRN = Chemical Abstracts Service Registry Number (if applicable); μg/L = micrograms per liter; pci/L = picocuries per liter; '-' = not applicable. \*Initial and Secondary COPC screening criteria are described in Section 5.3.







#### **6.1.2** Tentatively Identified Compounds

Tentatively Identified Compounds (TICs) were assessed where analyzed for both production and monitoring well samples. The Process Memo 97-005 User Guide (Coalition for Environmental Protection Restoration and Development [CEPRD], 2020) provides a description of TIC analysis:

"A TIC analysis uses the "library" within an analytical device to look for the chemical signature of 10,000s of chemicals. The purpose of a TIC analysis is to screen for the potential presence of chemicals that had not previously been identified. A separate analytical test method can then be used, if available, to confirm chemical identification and (for chemicals with analytical standards) to quantify the concentration of the chemical. Once verified and quantified, the water quality results can then be reviewed with DDW staff to determine next steps." (p .22)

The TICs results are tabulated and provided in Appendix C. A summary of those results is provided below.

- Reported TICs for production wells:
  - 13 analyses were completed for samples collected from SM-4, SM-8 and SM-9 in 2020.
  - Six detections were observed, of these detections:
    - three were reported as "Unknown" with concentrations ranging from 2.8 to 9.7 μg/L;
    - two were identified as PCE (EPA Method 625.1) with concentrations ranging from 12 to 13 μg/L
    - The remaining detection was identified as cis-2-octene with a reported concentration of 34 μg/L and was included in the raw water quality analysis.
- Reported TICs for monitoring wells:
  - 88 analyses were completed for samples taken from 13 locations in 2020.
  - 51 detections were observed, of these detections:
    - one was identified as 2(3H)-benzothiazolone with a reported concentration of 1.4 μg/L.
    - six were identified as cis-2-octene with concentrations ranging from 26 to 250 µg/L.
    - two were identified as benzothiazole with concentrations ranging from 4.0 to 8.5 μg/L.
    - one was identified as cyclohexyl isothiocyanate with a reported concentration of  $1 \mu g/L$ .
    - one was identified as 1-octene with a reported concentration of 25 μg/L.
    - five were identified as phthalic anhydride with concentrations ranging from 1.3 to 15 μg/L.
    - four were identified as PCE (EPA Method 625.1) with concentrations ranging from 12 to  $120 \mu g/L$ .
    - 38 were reported as "Unknown" with concentrations ranging 1.0 to 43 μg/L.

For TICs identified and reported with a Chemical Abstracts Service Registry Number (CASRN), statistical analysis was conducted and influent concentration estimates generated using the same approach adopted for all other constituents; the results of these data analyses are discussed in Section 6.1.3.







#### 6.1.3 **Microbiological Quality**

In accordance with the DDW Process Memo 97-005-R2020 the microbiological quality of the Olympic Well Field source water were evaluated and are summarized in this section. Groundwater samples collected<sup>3</sup> from the Olympic production and groundwater monitoring wells included the analyzes of the following microbiological parameters:

- Coliforms (total and fecal);
- Escherichia coli (E. coli); and
- Heterotrophic plate count (HPC).

The results are provided in Appendix D. The locations of the wells referenced below are shown in Figure 1-2.

Samples collected from Olympic production wells SM-4 and SM-9 were above the total coliforms' laboratory reporting limit of 1.1 MPN<sup>4</sup> /100ml at 6.9 and 2.2 MPN/100ml, respectively. Total coliforms were recorded at the reporting limit (1.1 MPN / 100 ml) in three groundwater monitoring wells (OB-3, OB-5 and SMB1-B#1).

All samples collected from the Olympic production wells were below the laboratory reporting limit for fecal coliforms (<1.1 MPN/100ml). However, there was one occurrence of fecal coliforms being reported above the laboratory reporting limit at 2.2 MPN /100 ml in the sample collected from groundwater monitoring well OB-15C.

HPC was reported above the laboratory reporting limit of 1 CFU<sup>5</sup>/ml in all samples collected with the exception of one groundwater monitoring well OB-12C. HPC was reported between 22 CFU /ml (SM-4) and 400 CFU /ml (SM-9). HPC ranged between 2 CFU /ml (OB-17B) and 400 CFU /ml (OB-11C) in samples collected from the monitoring wells (excluding OB-12C which was below the reporting limit).

E. coli was not detected in any of the samples collected.

#### 6.2 **Degradation of Chlorinated Solvents**

Chlorinated VOCs such as TCE and PCE are degradable under anaerobic conditions in groundwater; however, geochemical evidence (redox potential and dissolved oxygen) from groundwater samples in the Olympic Subbasin indicates that that the anaerobic conditions required for degradation of chlorinated VOCs either do not occur in groundwater in the subbasin, or occur only in limited settings.

The primary degradation pathway for PCE, TCE and other chlorinated VOCs is called dechlorination, a process by which the chlorine atoms are stripped from the ethylene molecule under strongly reducing (i.e., anaerobic or oxygen-free) conditions in groundwater. The degradation reaction is mainly driven by naturally-occurring soil microbes, and therefore is commonly called biodegradation. The microbes responsible for carrying out the degradation are strongly attached to aquifer soil particles and are not generally characterized in drinking water studies because they are not mobile in groundwater and are not

<sup>&</sup>lt;sup>3</sup>In accordance with the 97-005 Raw Water Characterization Sampling and Analysis Plan (Advisian 2020)

<sup>&</sup>lt;sup>4</sup> Most Probable Number







human pathogens. Microbiological characterization of drinking water quality is typically limited to coliforms, HPC, and E.Coli, which are characterized as part of the Step 2 Raw Water Quality Characterization (see Section 6.1.3) but are not associated with in-situ biodegradation of VOCs in groundwater.

The strongly reducing conditions needed for dechlorination reactions to occur fall within a sequence of oxidizing to reducing conditions that can occur in groundwater, which is illustrated in Figure 6-4. In the presence of dissolved oxygen, oxidizing conditions dominate in groundwater, and this is the most favorable condition for microbial growth and respiration. As oxygen is depleted in groundwater, other chemicals are used in place of oxygen to support microbial respiration, for example, nitrate (denitrifying conditions), ferric iron (iron reduction), manganese (manganese reduction), sulfate (sulfate reduction), and carbon dioxide (methane production or methanogenesis), in order of increasingly reducing conditions, and less favorable conditions energetically for the soil microbe population. In general, soil microbe populations will not use a less favorable chemical for respiration until all of the more favorable chemicals have been depleted. Commonly the population of soil microbes changes under different redox conditions.

As illustrated in Figure 6-4, dechlorination of TCE and PCE in groundwater occurs within the range of sulfate reduction and methane fermentation (Wiedemeier et al., 1999). Therefore, we do not expect to see this reaction occurring in groundwater that has significant concentrations of dissolved oxygen, nitrate, ferric iron, manganese (plus four valence state) or sulfate. Consequently, if dissolved oxygen or nitrate are present in in groundwater, it is unlikely that redox conditions are sufficiently anaerobic to support reductive dechlorination of TCE and PCE. The dechlorination of VOCs in groundwater produces a sequence of daughter products; i.e.: dechlorination of PCE or TCE results in the production of cis-1,2 DCE (and to a lesser extent, trans-1,2 DCE), and then vinyl chloride (VC), and finally ethene. The presence of these daughter products in groundwater is usually an indication that degradation of VOCs may be occurring in situ in the vicinity of the well, or in groundwater upgradient of the well.

Dissolved oxygen concentrations are typically measured in monitoring well samples during purging. In the B and C zones dissolved oxygen concentrations are variable, but average approximately 1 mg/L, representing aerobic conditions. Nitrate concentrations are typically in the range of 4.5 to 7.5 mg/L (as nitrogen [N]) in production wells SM-3 and SM-4; whereas for SM-8 and SM-9 which are screened deeper, nitrate concentrations are lower, 0.63 to 3.3 mg/L (as N) as measured in 2020. Cis-1,2 DCE concentrations are generally not detected in production well samples; however, trans-1,2-DCE has occasionally been detected in SM-4 at concentrations of 2 to 3 µg/L, suggesting that some degree of dechlorination of VOCs may be occurring within the capture zone of this well. VC has not been detected in the Olympic production wells, suggesting that if dechlorination is occurring within the capture zone, it is not strong enough to produce this penultimate daughter product. Moreover, the widespread occurrence of TCE at concentrations above the MCL in Olympic production wells strongly suggests that chlorinated VOC reduction is unlikely to be widespread in Olympic Subbasin groundwater, because redox conditions are not sufficiently reducing to support reductive dechlorination of VOCs. Consequently, natural attenuation of TCE and PCE concentrations by biodegradation is unlikely to be a significant process for reducing VOC concentrations in groundwater in the Olympic Subbasin.

The implications of this result for Step 2 RWQC characterization and treatment plant design are that the concentrations of VOCs in groundwater characterized in this Step 2 report are likely to be conservative, because there is no expectation that dechlorination will reduce VOC concentrations in the future influent

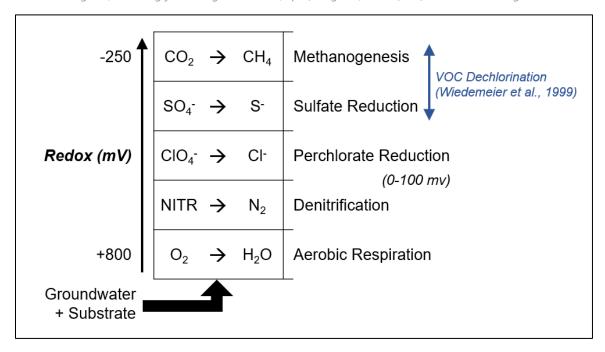






to the treatment plant, and therefore the treatment plant will be designed to manage VOC concentrations without relying on the reduction of VOC concentrations by biodegradation in situ.

Figure 6-1 Sequence of redox-driven degradation processes in groundwater, from aerobic conditions (bottom of diagram) to strongly reducing conditions (top of diagram). Modified from IRTC 2002 Figure 10-4.



# **6.3 Estimating Future Treatment Plant Influent Concentrations**

The results of the various calculation methods outlined in Section 5.4.1 used to derive estimates of future treatment plant influent concentrations are presented in the subsections below for the COPCs that meet both initial and secondary screening criteria (presented in Section 6.1). The results for all constituents within the groundwater quality data set are provided in Appendix E and Appendix F.

### **6.3.1** Statistical Analysis

A summary of the statistical analysis of the monitoring well groundwater quality data set is provided in Table 6-2. Statistical analysis results for all constituents within the monitoring well groundwater quality data set are provided in Appendix E.







Table 6-2 Summary of Statistical Analysis of Groundwater Monitoring Well Data for Each Capture Zone and Aquifer Zone for Key Synthetic Organic COPCs

			Concentrat	ion within (	Canturo	Concentrati	ion within Ca	nture Zone
Constituent	MCL	Aquifer	Zone Using U				an for Monit	
[Units]	or	Zone		/ell Data*	onitoring	Using Mea	Data**	ornig wen
[Ollits]	NL	Zone	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
1,1,2-		B-Zone	ND	-	ND	ND	-	ND
Trichloroethane		C-Zone	1.6	ND	ND	1.1	ND	ND
(1,1,2-TCA)	5	D-Zone	ND	ND	ND	ND	ND	ND
[µg/L]		Sunnyside	ND	ND	ND	ND ND	ND ND	ND
1,1-		B-Zone	0.67	-	0.15	0.52	-	0.15
Dichloroethane		C-Zone	0.40	0.90	0.13	0.243	0.60	0.13
(1,1-DCA)	5	D-Zone	ND	ND	ND	ND	ND	ND
[µg/L]		Sunnyside	ND	ND	ND	ND	ND	ND
1,1-		B-Zone	2.2	-	2.1	1.5	IND	1.7
Dichloroethene		C-Zone	2.0	3.9	0.22	1.5	3.3	0.22
(1,1-DCE)	6	D-Zone	0.5	0.5	0.22	0.5	0.5	0.22
(1,1-DCE) [μg/L]		Sunnyside		ND		ND		ND
			ND		ND		ND	
1,2,3-		B-Zone C-Zone	ND <b>0.0051</b>	0.016	ND ND	ND 0.0049	0.012	ND ND
Trichloropropane (1,2,3-TCP)	0.005		0.0031			0.0048	0.012 0.19	0.19
(1,2,3-1CP) [μg/L]		D-Zone	0.19	<b>0.19</b> 0.0041	<b>0.19</b> 0.0041	<b>0.19</b> 0.0041	0.19	0.0041
		Sunnyside		0.0041			0.0041	
1,2-		B-Zone	ND 0.33	0.50	ND	ND 0.33	0.47	ND
Dichloroethane	0.5	C-Zone	0.33	0.59	ND 0.15	0.33	0.47	ND 0.15
(1,2-DCA) [µg/L]		D-Zone	0.15	0.15	0.15	0.15	0.15	0.15
[µg/L]		Sunnyside	ND	ND	ND	ND FC 3	ND	ND 106.6
1.4.5:		B-Zone	102.1	-	117.4	56.3	-	106.6
1,4-Dioxane	1	C-Zone	27.8	48.4	1.8	22.7	36.3	2.3
[µg/L]		D-Zone	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND	ND	ND	ND
_		B-Zone	ND	-	ND	ND	-	ND
Benzene	1	C-Zone	0.23	1.6	ND	0.23	1.2	ND
[µg/L]		D-Zone	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND	ND	ND	ND
Carbon		B-Zone	0.63	-	ND	0.51	-	ND
Tetrachloride	0.5	C-Zone	0.68	0.62	0.12	0.57	0.52	0.12
(CTC)		D-Zone	0.3	0.3	0.3	0.3	0.3	0.3
[µg/L]		Sunnyside	ND	ND	ND	ND	ND	ND
cis-1,2-		B-Zone	0.26	-	ND	0.26	-	ND
Dichloroethene	6	C-Zone	0.85	49.4	0.8	0.74	30.4	0.61
(cis-1,2-DCE)		D-Zone	ND	ND	ND	ND	ND	ND
[µg/L]		Sunnyside	ND	ND	ND	ND	ND	ND
Methyl tert-butyl		B-Zone	0.53	-	ND	0.53	-	ND
ether (MTBE)	13	C-Zone	ND	3.8	ND	ND	1.8	ND
[µg/L]	, ,	D-Zone	ND	ND	ND	ND	ND	ND
[M3/ -]		Sunnyside	ND	ND	ND	ND	ND	ND
	5.1	B-Zone	3.1	-	2	1.9	-	2







Constituent [Units]	MCL or NL	Aquifer Zone	Concentrati Zone Using U W				ion within Ca an for Monito Data**	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
DEOA		C-Zone	0.97	0.89	1.4	0.67	0.66	1.4
PFOA		D-Zone	ND	ND	ND	ND	ND	ND
[ng/L]		Sunnyside	NA	NA	NA	NA	NA	NA
<del>-</del>		B-Zone	21.6	-	48.5	16.7	-	40.3
Tetrachloroethene	-	C-Zone	126.8	17.8	11.5	65.3	7.9	7.8
(PCE)	5	D-Zone	ND	ND	ND	ND	ND	ND
[µg/L]		Sunnyside	ND	ND	ND	ND	ND	ND
trans-1,2-		B-Zone	ND	-	ND	ND	-	ND
Dichloroethene	10	C-Zone	ND	0.97	0.63	ND	0.61	0.57
(trans-1,2-DCE)	10	D-Zone	ND	ND	ND	ND	ND	ND
[µg/L]		Sunnyside	ND	ND	ND	ND	ND	ND
T		B-Zone	2.2	-	7.5	1.7	-	6.6
Trichloroethene	-	C-Zone	126.9	19.2	3.8	63.6	13.0	3.1
(TCE)	5	D-Zone	1.5	1.5	1.5	1.5	1.5	1.5
[µg/L]		Sunnyside	ND	ND	ND	ND	ND	ND
		B-Zone	ND	-	ND	ND	-	ND
Vinyl Chloride	0.5	C-Zone	ND	1.9	ND	ND	1.3	ND
[µg/L]	0.5	D-Zone	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND	ND	ND	ND

#### Notes:

μg/L = micrograms per liter; ng/L = nanograms per liter; 'ND' = non-detect; 'NA' = not available; '-' = not applicable (production well is not screened in aquifer zone). Values above respective MCLs or NLs are highlighted **bold**.

\*UCL95 (95 percent upper confidence limit of the population mean) based on maximum suggested UCL95 statistics produced by ProUCL software (EPA 2015) calculated for censored data sets (using all detects and non-detects).

\*\*Mean calculated for censored data sets (using all detects and non-detects) via Kaplan Meier Method (using ProUCL software; EPA 2015).

In the event either the mean or UCL95 statistics could not be calculated, the maximum value was adopted as a conservative approach. For non-detect (ND) results a value of 0 was assumed; where data was not available (NA) for a constituent in any given layer and/or capture zone, a value of 0 was assumed.

For the D-Zone, monitoring well SMB-1-B#1 was used for all production well capture zones; for the Sunnyside, monitoring well SMB-1 #4 was used for all production well capture zones.

#### 6.3.2 Production Wells Concentration Estimates Results

The results of the evaluation of estimated future concentrations for all constituents within the three Olympic Well Field groundwater production wells (i.e., SM-4, SM-8 and SM-9) are provided in this section.

Table 6-3 summarizes flow-weighted calculation results which are the first component of the analysis. This includes transmissivities for production well screens, calculated by multiplying total screen length in each aquifer zone by the hydraulic conductivity for the applicable layer from the calibrated groundwater model. The transmissivities across each well screen was then used to calculate the proportion of flow to the well from each aquifer zone, i.e., the proportion of total flow in the production well coming from each screen interval.







Table 6-4 provides the flow-weighted concentration estimates for each production well applying the flow calculations in Table 6-3 to the mean and UCL95 statistics for each of the COPCs that meet both initial and secondary screening criteria (presented in Section 6.1). Production wells concentration estimates for all constituents within the groundwater quality data set are provided in Appendix E.

Table 6-3 Flow-Weighting Calculations for Production Wells

Aquifer Zone (Model		al Pum <sub>l</sub> ite (gpi		Lengt	tal Scre th in Ac Cone (ft	quifer	Con	lydraul ductivi er Zone	ty in	Aqı	smissiv uifer Zo (ft²/d)	one		ated Fl uifer Zo (gpm)	-
Layer)	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
B-Zone				25	0	25	16.52	26.97	1.16	413	0	29	211	0	11
C-Zone	700	700	600	75	55	80	3.06	1.29	1.19	230	71	95	117	30	35
D-Zone	700	700	600	115	30	40	2.17	3.04	2.28	250	91	91	127	39	34
Sunnyside				45	110	170	10.65	13.44	8.27	479	1,478	1,406	245	631	520

**Notes:**  $gpm = gallons \ per \ minute$ ; ft = feet;  $ft/d = feet \ per \ day$ ;  $ft2/d = square \ feet \ per \ day$ .

Table 6-4 Estimated Flow-Weighted Concentrations in Individual Production Wells for Key Synthetic Organic COPCs

Constituent [Units]	MCL or NL		ed Conce uction W UCL95*	ntrations ells Using		ed Conce uction W Mean**	ells Using	Concen	iction V trations 020***	
		SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
1,1,2-Trichloroethane [µg/L]	5	0.27	ND	ND	0.18	ND	ND	ND	ND	ND
1,1-Dichloroethane (1,1-DCA) [µg/L]	5	0.27	0.039	0.01	0.2	0.026	0.01	0.16	0.12	ND
1,1-Dichloroethene (1,1-DCE) [µg/L]	6	1.1	0.2	0.079	0.81	0.17	0.071	1.2	0.72	ND
1,2,3- Trichloropropane (1,2,3-TCP) [µg/L]	0.005	0.037	0.015	0.014	0.037	0.015	0.014	0.0026	ND	ND
1,2-Dichloroethane [µg/L]			0.034	0.0085	0.083	0.029	0.0085	ND	ND	ND
1,4-Dioxane [µg/L]	1	35.4	2.1	2.2	20.8	1.6	2	20	5.4	3.3
Benzene [µg/L]	1	0.039	0.068	ND	0.039	0.051	ND	ND	ND	ND







Constituent [Units]	MCL or NL		ed Conce uction W UCL95*	ntrations ells Using		ed Conce uction W Mean**	ntrations ells Using	Concen	uction V trations 020***	
		SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
Carbon Tetrachloride (CTC) [µg/L]	0.5	0.36	0.044	0.024	0.3	0.039	0.024	0.13	0.15	ND
cis-1,2- Dichloroethene (cis-1,2-DCE) [μg/L]	6	0.22	2.1	0.047	0.2	1.3	0.036	0.45	0.22	ND
Methyl tert-butyl ether (MTBE) [µg/L]	13	0.16	0.17	ND	0.16	0.077	ND	ND	ND	ND
PFOA [ng/L]	5.1	1.1	0.039	0.12	0.68	0.028	0.12	0.96	ND	2.9
Tetrachloroethene (PCE) [µg/L]	5	27.7	0.77	1.5	16	0.34	1.2	54	0.39	41
trans-1,2- Dichloroethene [µg/L]	10	ND	0.042	0.037	ND	0.026	0.033	ND	ND	ND
Trichloroethene (TCE) [µg/L]	5	22.2	0.92	0.44	11.4	0.65	0.38	59	7	0.34
Vinyl Chloride [μg/L]	0.5	ND	0.081	ND	ND	0.058	ND	ND	ND	ND

#### Notes:

 $\mu g/L = micrograms\ per\ liter;\ ng/L = nanograms\ per\ liter;\ 'ND' = non-detect;\ '-' = not\ applicable.$  Values above respective MCLs or NLs are highlighted **bold**.

#### **6.3.3** Treatment Plant Influent Concentration Estimates

Estimated future treatment plant influent concentrations (comprising the combined flow from the three production wells) using estimates derived from UCL95 statistics for each of the COPCs that met both initial and secondary screening criteria (presented in Section 6.1) are provided in Table 6-5. In addition, estimates with a safety factor of 1.5 (1.2 for 1,2,3-TCP) applied to constituent concentration are also provided in Table 6-5. Finally, for information, influent estimates using the maximum observed value from recent production well sampling data collected during 2020 is also provided in Table 6-5. Treatment plant influent concentration estimates for all constituents within the groundwater quality data set are provided in Appendix F.

Results indicate four COPCs are projected to be at concentrations above their respective MCL or NL in treatment plant influent, including 1,4-dioxane, PCE, TCE and 1,2,3-TCP. For comparison, influent estimates

<sup>\*</sup>UCL95 from statistical analysis of the monitoring well groundwater quality data set (described in Section 6.3.1).

<sup>\*\*</sup>Mean from statistical analysis of the monitoring well groundwater quality data set (described in Section 6.3.1).

<sup>\*\*\*</sup>Maximum observed values from production wells from available 2020 sampling data (provided for information only). In the event either the mean or UCL95 statistics could not be calculated, the maximum value was adopted as a conservative approach.







based on recent (2020) water quality data from production wells indicate that 1,4-dioxane, PCE and TCE would be currently above their respective MCL or NL, with 1,2,3-TCP being below its MCL based on these data (as shown in Table 6-5).

Table 6-5 Estimated Concentrations in Treatment Plant Influent for Key Synthetic Organic COPCs

		Using U	CL95*	Using Produ Concentrations	
Constituent [Units]	MCL or NL	Plant Influent Concentration Estimates	With Safety Factor Applied***	Plant Influent Concentration Estimates	With Safety Factor Applied***
1,1-Dichloroethane (1,1-DCA) [μg/L]	5	0.11	0.17	0.098	0.15
1,1,2-Trichloroethane (1,1,2-TCA) [μg/L]	5	0.094	0.14	ND	ND
1,1-Dichloroethene (1,1-DCE) [μg/L]	6	0.47	0.71	0.67	1.0
1,2,3-Trichloropropane (1,2,3- TCP) [µg/L]	0.005	0.022	0.026	0.00091	0.0011
1,2-Dichloroethane (1,2-DCA) [µg/L]	0.5	0.043	0.065	ND	ND
1,4-Dioxane [μg/L]	1	13.8	20.7	9.9	14.9
Benzene [μg/L]	1	0.037	0.056	ND	ND
Carbon Tetrachloride (CTC) [µg/L]	0.5	0.15	0.23	0.098	0.15
cis-1,2-Dichloroethene (cis-1,2-DCE) [µg/L]	6	0.84	1.26	0.23	0.35
Methyl tert-butyl ether (MTBE) [µg/L]	13	0.11	0.17	ND	ND
PFOA [ng/L]	5.1	0.44	0.66	1.2	1.8
Tetrachloroethene (PCE) [µg/L]	5	10.4	15.6	31	46.5
trans-1,2-Dichloroethene (trans-1,2-DCE) [µg/L]	10	0.026	0.039	ND	ND
Trichloroethene (TCE) [μg/L]	5	8.2	12.3	23	34.5
Vinyl Chloride [μg/L]	0.5	0.028	0.042	ND	ND

**Notes:**  $\mu g/L = micrograms \ per \ liter; \ ng/L = nanograms \ per \ liter; \ 'ND' = non-detect.$ 

Values above respective MCLs or NLs are highlighted **bold**.

<sup>\*</sup> From production wells concentration estimates (described in Section 6.3.2).

<sup>\*\*</sup>Maximum observed values from production wells from available 2020 sampling data (provided for information only).







\*\*\*A safety factor of 1.5 was applied for each constituent, except for 1,2,3-TCP which used a safety factor of 1.2.

#### **6.3.4** Uncertainty and Safety Factors

In accordance with the DDW Process Memo 97-005 -R2020 (DDW 2020), the degree of uncertainty associated with the estimated influent concentrations was evaluated, and safety factors commensurate with the degree of uncertainty provided. These requirements are addressed below.

Data limitations and uncertainties include, but are not limited to the following:

- Locations within projected capture zones for production wells that were not sampled may have different water quality from those areas where sampling data is available from established wells. It is noted that no monitoring network is robust enough to describe all subsurface conditions; hence, unsampled locations may differ from those interpreted from established wells.
- The potential for unknown (i.e., yet unidentified) historical and future contamination source(s) and pathway conditions in the vicinity of the production wells.
- Uncertainties in historical and future well field operations, including changes from modeled pumping due to operational demands and/or regulatory changes.

To account for such uncertainties, comprehensive data analyses were carried out, as presented herein, including:

- Spatial and statistical analysis of monitoring well data to characterize the distribution and persistence
  all constituents in groundwater within the projected capture zones and to provide an indication of the
  quality of water is anticipated to arrive at the production wells in the future, including provision of a
  range of potentially likely and conservative statistics and future production well concentration
  estimates using mean and UCL95 statistics.
- Trend and correlation analysis of production well data to characterize constituents in historical raw water quality at the production wells.

With respect to treatment design and safety factors, the following were considered:

- Conservative influent water quality assumptions adopted for calculations (as discussed above).
  - Future concentration estimates are considered conservative as they assume there are no transport mechanisms that reduce constituent concentrations, such as attenuation, retardation or degradation.
  - Future concentration estimates assume all flow from each aquifer zone into each production well screen has concentrations equal to the adopted value from monitoring well statistics (mean or UCL95 statistics, or maximum observed value if either of these statistics were unavailable). This may result in an over-estimate for production well influent concentrations for some constituents as there are expected to be areas within each capture zone which have low or non-detect concentrations of the constituents evaluated.
  - For treatment plant influent estimates, only the production well estimates derived from UCL95 statistics were used as this is more conservative than the mean.
- A safety factor 1.5 (1.2 for 1,2,3-TCP) was adopted for this RWQC for influent concentrations which spans the range of potential concentrations developed herein and described above. Average water







quality conditions are less than the design conditions which will include a safety factor and therefore provide an additional safety measure for mitigating unexpected COPC concentrations.

#### 6.4 Trend Analysis

This section presents the results of the Mann-Kendall and Sen's Slope trend analyses for water quality data from production wells SM-3 and SM-4 using the approach outlined in Section 5.4.2. The tabulated results of the Mann-Kendall and Sen's Slope trend analysis are provided in Appendix G, and trend charts generated for each COPC for visual trend assessment purposes are provided in Appendix H. A summary of the results is presented below.

For production well SM-3, statistically significant and visual decreasing trends were identified for 1,4-dioxane when assessing the data from 2014 to  $2020^6$  and PCE when assessing data from 2012 to 2020. In contrast, data from 2017 to 2020 did not contain any statistically significant trends. Visually, nitrate and nitrite as nitrogen (as N) has exhibited variability since 2016. Other COPCs with concentrations above the detection limit, including TCE, were deemed stable. 1,2,3-TCP concentrations were below the detection limit of 0.005  $\mu$ g/L in samples collected since its incorporation in the quarterly Olympic Well Field monitoring analytical schedule in 2018.

For production well SM-4, statistically significant and visual increasing trends were identified for PCE when using both the 2012 to 2020 and 2017 to 2020 data sets. A statistically significant and visual decreasing trend was identified for 1,4-dioxane in the 2017 to 2020 data. Similar to production well SM-3, visual assessment indicated nitrate and nitrite (as N) has exhibited variability since 2016. Higher concentrations of TCE were recorded between 2012 and 2014, with an observable rebound to similar levels noted in late 2018 to 2020; however, no statistically significant trend was identified. Visual assessment of 1,2,3-TCP indicated a decreasing trend in 2018; however, since 2019 concentrations have been generally stable. No statistically significant trend was identified for 1,2,3-TCP. Visual assessment of other COPCs generally indicated that parameters with concentrations above detection limit have remained stable during the period from 2012 to 2020.

## 6.5 Raw Water Quality Variability

This section presents the results of COPC concentration variability with pumping rate and time (seasonal and long-term) observed in the production wells SM-3 and SM-4 using the approach outlined in Section 5.4.3.

#### 6.5.1 Variation with Pumping Rate

This section addresses observed relationships between production (pumping) at a single production well and the concentration of each COPC. Concentration data and monthly pumping volumes from January 1, 2012 to June 30, 2020 are shown in the charts provided in Appendix H. The charts were used to assess potential concentration changes with respect to pumping.

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<sup>&</sup>lt;sup>6</sup> 1,4-dioxane testing commenced in 2014 when it was identified as an emerging contaminant of concern, hence the full 1,4-dioxane data series is 2014-2020







It is important to note that Olympic Well Field production well SM-3 was taken out-of-service in May 2013 due to well casing failure. During the shutdown, the well was re-lined and brought back into production in August 2014. The diameter of the screen and casing reduced as a result of re-lining which in-turn resulted in reduced production rates at SM-3. These reduced production rates are shown in the charts generated for SM-3 in Appendix H. SM-4 has generally been pumping consistently since 2012, with the exception of a short-duration decrease in rate in 2013, an overall decrease in rate since late 2016, and low to no pumping in early 2020.

As noted previously in the approach outlined in Section 5.4.3, the data set was visually assessed to determine correlations between COPC concentrations and pumping, with a positive correlation being defined as instances where the recorded concentration of a parameter increases (or decreases) as pumping rate increases (or decreases). On the other hand, a negative correlation is defined as instances where the recorded concentration of a parameter increases (or decreases) as the pumping rate decreases (or increases).

Based on visual assessment, a positive correlation with pumping rate was noted for the following COPCs for SM-3:

- 1,4-dioxane concentrations were higher (approximately 4 to 8 μg/L) in 2012-2013 when pumping rates were greater. Once pumping rates decreased in 2014, 1,4-dioxane concentrations remained below 2 μg/L.
- Nitrate concentrations were higher in the 2012-2013 when production was greater.
- TCE concentrations decreased when pumping ceased in 2013 and remained lower when reduced pumping resumed in 2014. Additional decreases were noted following short periods of no pumping in 2015, 2018, and 2019. This positive correlation between SM-3 pumping and TCE concentration can be seen in the chart presented in Based on visual assessment, a negative correlation with pumping rate was noted at SM-4 for the following COPCs:
- An upward trend identified for 1,4-dioxane from 2017 to 2020 may be associated with decreasing pumping rates

With reference to Figure 6-3 below, a negative correlation between PCE concentration and pumping is observed since 2017.

No clear correlation with pumping rate was noted for TCE and 1,2,3-TCP for SM-4.







- Figure 6-2 below.
- No clear correlation with pumping rate was noted for PCE
- 1,2,3-TCP concentrations were below the detection limit in all samples collected from SM-3.

Based on visual assessment, a negative correlation with pumping rate was noted at SM-4 for the following COPCs:

• An upward trend identified for 1,4-dioxane from 2017 to 2020 may be associated with decreasing pumping rates

With reference to Figure 6-3 below, a negative correlation between PCE concentration and pumping is observed since 2017.

No clear correlation with pumping rate was noted for TCE and 1,2,3-TCP for SM-4.







Figure 6-2 Production Well SM-3 - TCE Concentration versus Pumping Chart (from Appendix H)

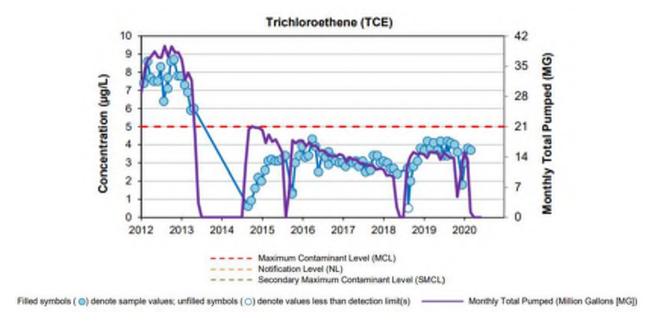
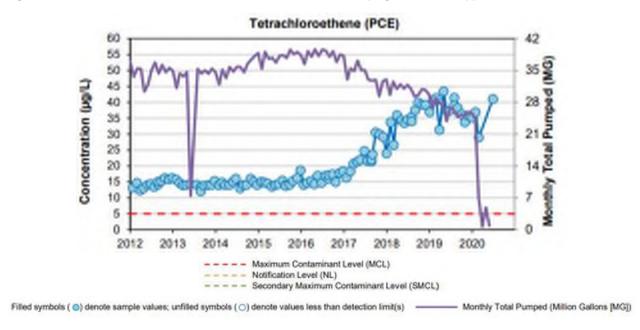


Figure 6-3 Production Well SM-4 - PCE Concentration versus Pumping Chart (from Appendix H)



With reference to the charts provided in Appendix H and SM-3, concentrations of COPCs in the Olympic Well Field, e.g., 1,4-dioxane, TCE and PCE, reduced in response to decreased production from the well since 2014. It is conjectured that reduced (~50%) pumping at SM-3 resulted in less contamination capture, with the higher pumping at well SM-4 drawing contamination in, and away from SM-3.







Also, of interest, is the observed increase in concentrations of COPCs in response to reduced production at SM-4. Since the well is located proximal to the two key contamination source sites (refer to Section 3) and therefore close to the core of the contaminant plume(s), it is possible that decreased production resulted in a reduced capture zone, which in-turn resulted in less "uncontaminated" groundwater capture by SM-4, which subsequently reduced the effects of in-well dilution.

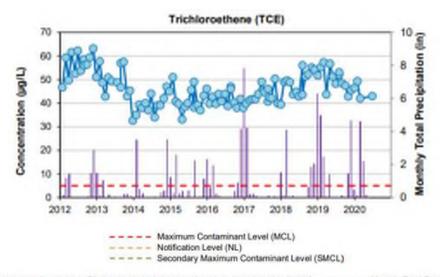
#### 6.5.2 Variations with Time (Seasonal Changes and Wet / Dry Periods)

Seasonal impacts on COPC concentrations in the temporal production well data sets were assessed as part of this RWQC study. The objective of this evaluation was to determine whether changes in concentrations of COPCs could be correlated to seasonal variation in precipitation. Variability of the climate over time was also assessed to determine whether changes in concentrations of COPCs could be correlated to periods considered to be "wet" (i.e., periods of high precipitation) versus "dry" (i.e., periods of low precipitation), as outlined previously in Section 5.4.3.

Over the 2012 to 2020 observation period, climatic conditions became consistently wetter (see charts in Appendix I). Therefore, the greatest contrast between wet and dry periods occurred between the earliest part of the observation period to the most recent portion of the data set from 2020.

Based on a visual assessment of the data, no correlation with monthly precipitation and wet/dry periods were observed for any parameters at SM-3 and SM-4. For example, Figure 6-4 below presents temporal (2012-2020) TCE concentration and precipitation data. No discernable correlation between seasonal changes or wet/dry periods and TCE concentrations is evident.

Figure 6-4 Production Well SM-4 - TCE Concentration versus Precipitation Chart (from Appendix I)



Filled symbols ( (ii)) denote sample values; unfilled symbols ( (iii)) denote values less than detection limit(s) ———— Monthly Total Precipitation (inches (in))







# 7 Conclusions

The DDW 97-005 process consists of 10 steps for evaluating the proposed use of an extremely impaired source for direct potable use. A source is considered to be "extremely impaired" if it meets two or more of 10 DDW-developed criteria. Based on the evaluation of available water quality data, groundwater in the vicinity of the Olympic Well Field meets up to three criteria. The Olympic Well Field RWQC study (*this report*) is Step 2 of the 10-step process and satisfies the applicable requirement of the DDW Process Memo 97-005-R2020. A summary of the key outcomes of this RWQC study is provided below.

A series of data analyses were conducted to evaluate raw water quality to meet the primary objective of characterizing influent water quality that will enter the planned treatment system, so that an appropriate level of monitoring and treatment can be designed. This included screening of water quality data against specific criteria and regulatory values to identify COPCs, analyzing water quality data to estimate future treatment plant influent concentrations, trend analysis for historical water quality to understand how water quality trends have changed over time, and analysis of variability to develop an understanding of how water quality has changed under the influence of certain factors such as pumping and seasonal variation in precipitation.

An initial screening of water quality data indicated 42 COPCs based on constituents detected in production wells (or constituents with three or more detections in monitoring wells), with a ratio of maximum concentration to MCL or NL greater than 0.05 (5%). The list of COPCs was then further refined to determine which would drive treatment system design (using UV/AOP and GAC) based on constituents which are synthetic organics and had a ratio of maximum concentration to MCL or NL greater than 0.5 (50%). This yielded a list of 15 synthetic organic COPCs, as follows:

- Nine COPCs from production well data:
  - 1,1-dichloroethane (1,1-DCA)
  - 1,1-dichloroethene (1,1-DCE)
  - 1,2,3-trichloropropane (1,2,3-TCP)
  - 1,4-dioxane
  - carbon tetrachloride
  - cis-1,2-dichloroethene (cis-1,2-DCE)
  - perfluorooctanoic acid (PFOA)
  - tetrachloroethene (PCE)
  - trichloroethene (TCE)

- An additional six COPCs from monitoring well data:
  - 1,1,2-trichloroethane (1,1,2-TCA)
  - 1,2-dichloroethane (1,1-DCA)
  - benzene
  - methyl tert-butyl ether (MTBE)
  - trans-1,2-dichloroethene (trans-1,2-DCE)
  - vinyl chloride

Statistical analysis and flow-weighting calculations were then conducted to estimate future treatment plant influent concentrations using monitoring well groundwater quality data for all constituents including the identified COPCs. The results of this analysis indicated four COPCs are projected to be at concentrations above their respective MCL or NL in treatment plant influent, including 1,4-dioxane, PCE, TCE and 1,2,3-TCP, as summarized in Table 7-1. For comparison, influent estimates based on recent (2020) water quality data from production wells indicate that 1,4-dioxane, PCE and TCE would be currently above their respective MCL or NL, with 1,2,3-TCP being below its MCL based on this data, as shown in Table 7-1.







Table 7-1 Summary of Estimated Concentrations in Treatment Plant Influent for Key Synthetic Organic COPCs

		Using UC	L95*	Using Production We From 202	
Constituent [Units]	MCL or NL	Plant Influent Concentration Estimates	With Safety Factor Applied***	Plant Influent Concentration Estimates	With Safety Factor Applied***
1,2,3-TCP [μg/L]	0.005	0.022	0.026	0.00091	0.0011
1,4-Dioxane [µg/L]	1	13.8	20.7	9.9	14.9
PCE [μg/L]	5	10.4	15.6	31	46.5
TCE [µg/L]	5	8.2	12.3	23	34.5

**Notes:**  $\mu g/L = micrograms per liter; values above respective MCLs or NLs are highlighted$ **bold**.

Trend analysis was performed on available temporal production well data for COPCs to inform potential future concentration trends and design of the treatment system. Key outcomes from the trend analysis include the identification of statistically significant upward trends for PCE and 1,4-dioxane. Although no statistically significant trend was identified for TCE, visual assessment of trend charts indicates that recent (2018-2020) concentrations are rebounding to elevated levels recorded between 2012 and 2014. No statistically significant or visual trend was identified for 1,2,3-TCP. Visual assessment indicated nitrate and nitrite (as nitrogen) exhibited variability since start of analysis period (2012) with no discernible trend identified. Visual assessment of other COPC trends generally indicated that parameters with concentrations above detection limit remained stable during the analysis period (2012-2020).

An assessment of concentration variability generally indicates that concentrations of 1,4-dioxane, TCE, PCE exhibit an observable correlation to operational status (i.e., pumping or non-pumping) and pumping rate of a well or at nearby wells; however, seasonal impacts or seasonal signals are not distinguishable in the data sets.

<sup>\*</sup> From production wells concentration estimates.

<sup>\*\*</sup>Maximum observed values from production wells from available 2020 sampling data (provided for information only).

<sup>\*\*\*</sup>A safety factor of 1.5 was applied for each constituent, except for 1,2,3-TCP which used a safety factor of 1.2.







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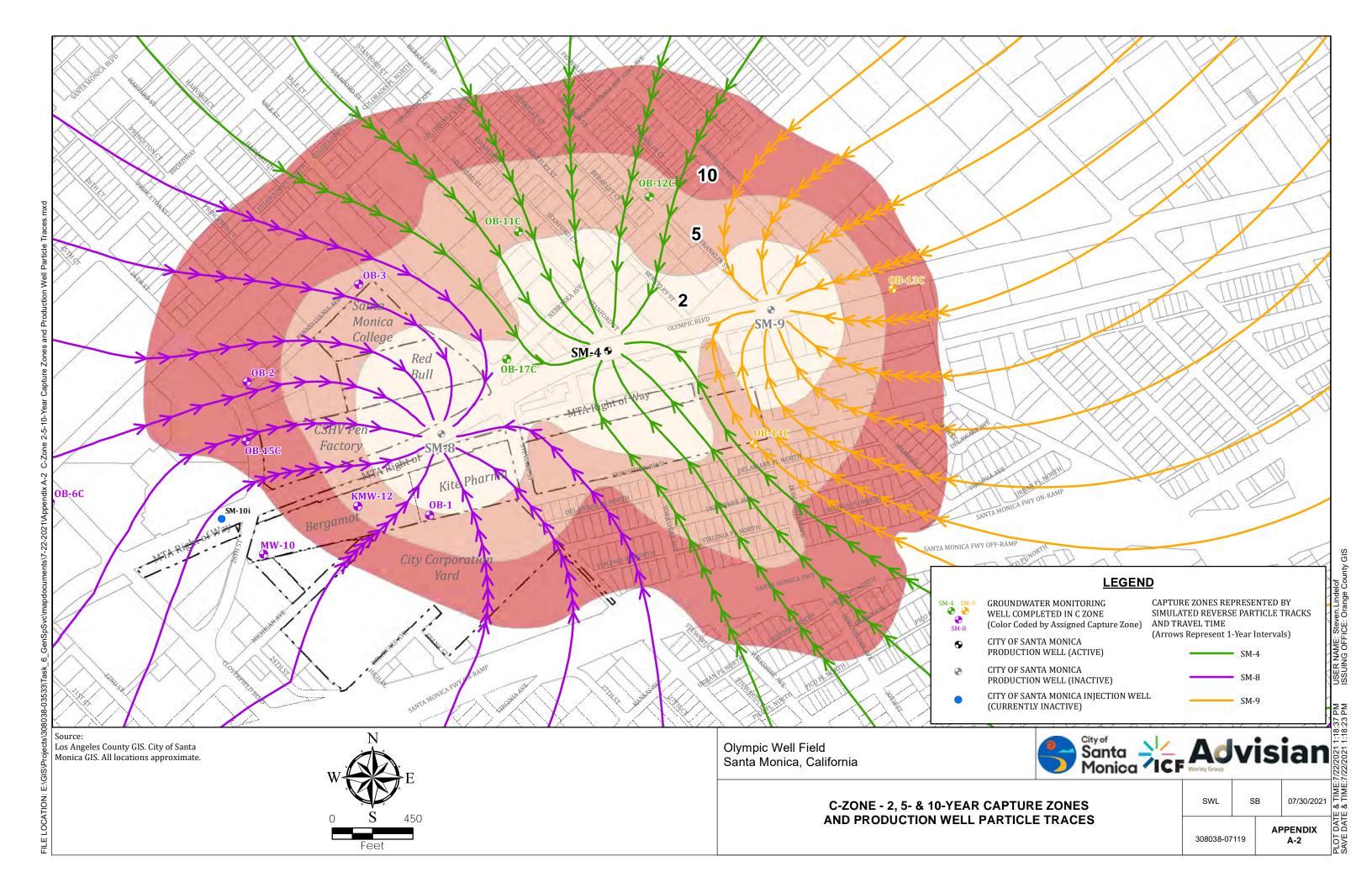


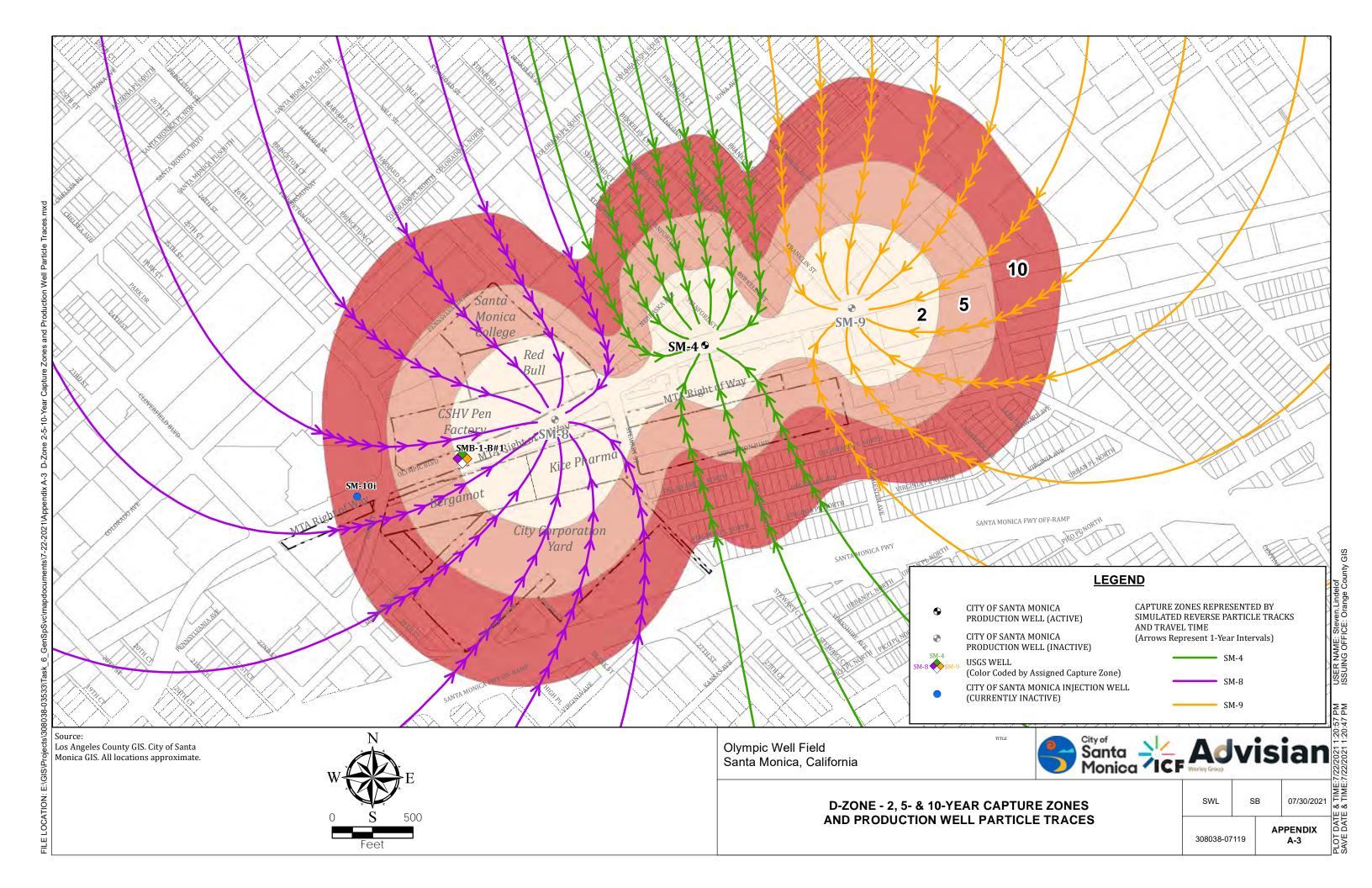


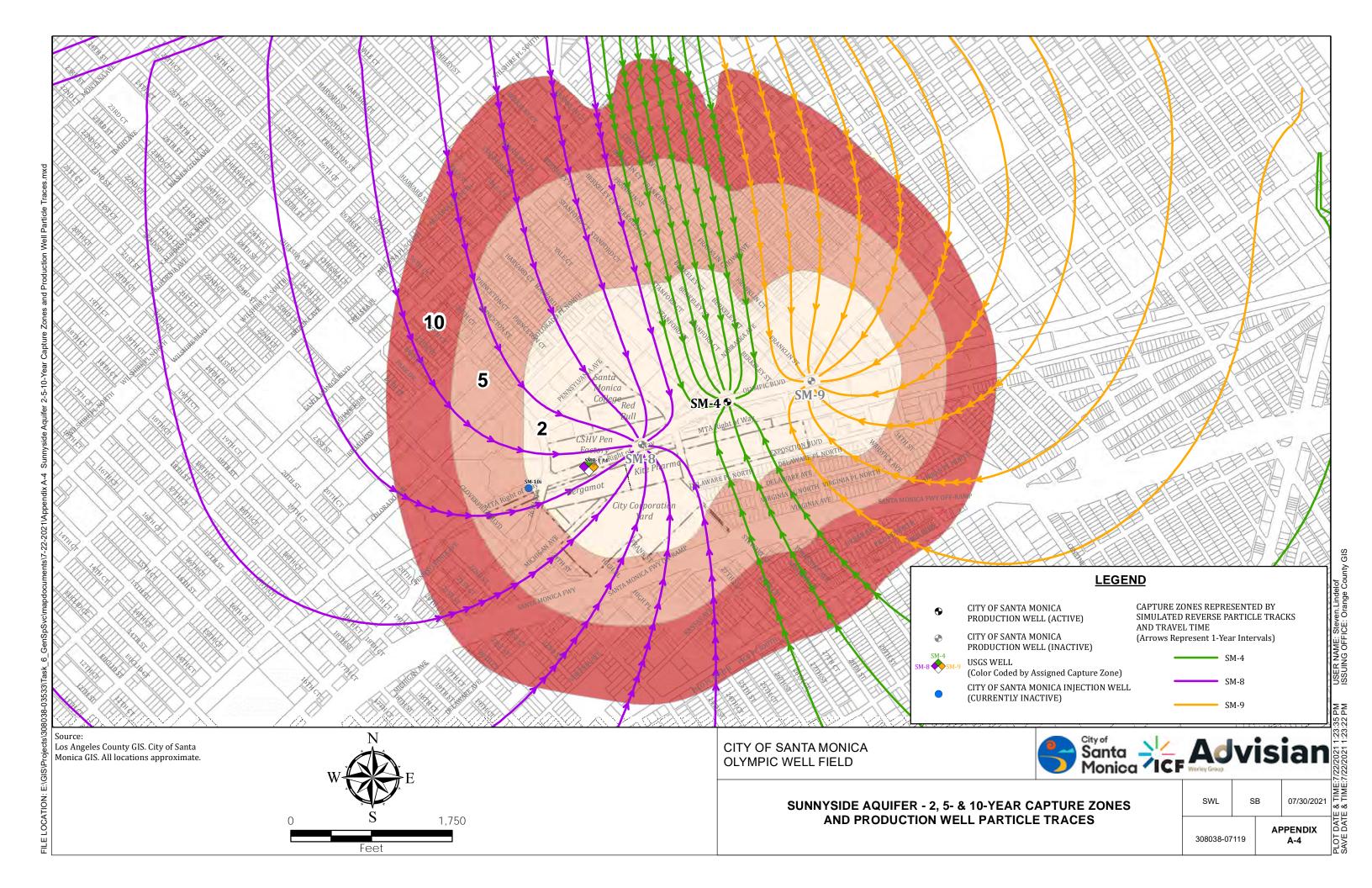


# Appendix A Modeled Capture Zones and Monitoring Well Locations















# **Appendix B Screening Evaluation Results**







						Producti	on Wells			Monitori	ng Wells			Mee	ets COPC Crite		ing
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Init	tial 3 4	Seco 5	ondary 6
3567-62-2	1-(3,4-Dichlorophenyl)-3-methylurea	μg/L	-	-	3	0	<1	-	13	0	<1	-					
2327-02-8	1-(3,4-Dichlorophenyl)urea	μg/L	-	-	3	0	<1	-	13	0	<1	-					
630-20-6	1,1,1,2-Tetrachloroethane	μg/L	-	-	34	0	<0.5	-	475	0	<1	-					
71-55-6	1,1,1-Trichloroethane	μg/L	200	-	201	0	<0.5	-	475	0	<1	-			✓		
79-34-5	1,1,2,2-Tetrachloroethane	μg/L	1	-	34	0	<0.5	-	475	0	<1	-			✓		
79-00-5	1,1,2-Trichloroethane	μg/L	5	-	34	0	<0.5	-	475	8	11	2.20		✓	✓ ✓	✓	✓
75-34-3	1,1-Dichloroethane	μg/L	5	-	201	2	0.16	0.03	475	67	7.1	1.42	✓	✓	✓ ✓	✓	✓
75-35-4	1,1-Dichloroethene	μg/L	6	-	201	117	2.9	0.48	475	138	32	5.33	✓	✓	✓ ✓	✓	✓
563-58-6	1,1-Dichloropropene	μg/L	-	-	34	0	<0.5	-	475	0	<1	-			1		
57-14-7	1,1-Dimethylhydrazine	μg/L	-	-	3	0	<2	-	13	0	<2	-					
87-61-6	1,2,3-Trichlorobenzene	μg/L	-	-	34	0	<0.5	-	475	0	<1	-					
96-18-4	1,2,3-Trichloropropane	μg/L	0.005	_	38	17	0.01	2.00	609	29	0.19	38.00	✓	✓	<b>✓ ✓</b>	✓	<b>✓</b>
120-82-1	1,2,4-Trichlorobenzene	μg/L	5	_	37	0	<1	-	494	0	<10	-			<b>✓</b>		
95-63-6	1,2,4-Trimethylbenzene	μg/L	_	330	34	0	<0.5	_	475	0	<1	_			<b>√</b>	+	+
96-12-8	1,2-Dibromo-3-chloropropane	μg/L	0.2	-	16	0	<0.01	_	475	0	<5	_			<b>✓</b>		
106-93-4	1,2-Dibromoethane	μg/L μg/L	0.05	_	16	0	<0.02	_	475	0	<del></del>	_			<b>✓</b>	_	
95-50-1	1,2-Dichlorobenzene	μg/L μg/L	600	-	37	0	<1	_	494	0	<10	_			<b>√</b>	_	
107-06-2	1,2-Dichloroethane	μg/L μg/L	0.5	_	201	0	<0.5	_	475	21	6.6	13.20		<b>√</b>	√ √	<b>√</b>	<b>✓</b>
78-87-5	1,2-Dichloropropane	μg/L μg/L	5	_	34	0	<0.5		475	1	0.16	0.03		•	√ ·	+	+
122-66-7	1,2-Dichloroproparie	μg/L μg/L	J	_	3	0	<1	_	19	0	<10	0.03				+	+
108-67-8	1,3,5-Trimethylbenzene	μg/L μg/L		330	34	0	<0.5		475	0	<1				<b>√</b>		
99-35-4	1,3,5-Trinitrobenzene		-	-	6	0	<10	-	32	0	<100						1
		μg/L	-		37	_		-	494	, and the second							1
541-73-1	1,3-Dichlorobenzene	μg/L	-	-	34	0	<1 <0.5	-	494	0	<10	-					4
142-28-9	1,3-Dichloropropane	μg/L	- 0.5	-		0		-		-	<1	-			<b>√</b>		4
542-75-6	1,3-Dichloropropene	μg/L	0.5	-	34	0	<0.5	-	19	0	<0.5	-	<b>✓</b>		<u> </u>		4
99-65-0	1,3-Dinitrobenzene	μg/L	-	-	3	ı	0.32	-	13	ı	0.26	-	<b>V</b>		<b>√</b>		4
106-46-7	1,4-Dichlorobenzene	μg/L	5	-	37	0	<1	-	494	0	<10	-					1
123-91-1	1,4-Dioxane	μg/L	-	1	68	64	26	26.00	413	367	440	440.00	<b>√</b>	<b>√</b>	✓ ✓	<b>√</b>	<b>Y</b>
763051-92-9	11CI-PF3OUdS	μg/L	-	-	8	2	0.0017	-	16	0	<0.002	-	✓				4
57-91-0	17-a-Estradiol	μg/L	-	-	3	0	<0.001	-	13	0	<0.001	-					4
57-63-6	17-a-Ethynylestradiol	μg/L	-	-	2	0	<0.001	-	13	0	<0.001	-					4
50-28-2	17-b-Estradiol	μg/L	-	-	3	0	<0.001	-	13	0	<0.03	-					4
124-19-6	1-Nonanal	μg/L	-	-	3	0	<2	-	14	0	<2	-			$\longrightarrow$		4
934-34-9	2(3H)-Benzothiazolone	μg/L	-	-	0	0	-	-	1	1	1.4	-					4
594-20-7	2,2-Dichloropropane	μg/L	-	-	34	0	<0.5	-	475	0	<1	-					4
51207-31-9	2,3,7,8-Tetrachlorodibenzofuran	μg/L	-	-	3	0	<0.00005	-	19	0	<0.00005	-					4
93-76-5	2,4,5-T	μg/L	-	-	5	0	<0.2	-	19	0	<0.2	-					
93-72-1	2,4,5-TP (Silvex)	μg/L	50	-	11	0	<1	-	19	0	<0.2	-			✓		4
95-95-4	2,4,5-Trichlorophenol	μg/L	-	-	3	0	<1	-	19	0	<10	-					4
88-06-2	2,4,6-Trichlorophenol	μg/L	-	-	3	0	<1	-	19	0	<10	-					
118-96-7	2,4,6-Trinitrotoluene	μg/L	-	1	3	0	<5	-	13	0	<5	-			✓		
94-75-7	2,4-D	μg/L	70	-	15	0	<10	-	19	0	<0.4	-			✓		
94-82-6	2,4-DB	μg/L	-	-	5	0	<2	-	19	0	<2	-					
120-83-2	2,4-Dichlorophenol	μg/L	-	-	3	0	<1	-	19	0	<10	-					
105-67-9	2,4-Dimethylphenol	μg/L	-	-	3	0	<1	-	19	0	<10	-					







						Producti	on Wells			Monitor	ing Wells		١	Meets	COPC	Screen ria*	ning
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Initio	al 3 4	Second 5	ondary 6
51-28-5	2,4-Dinitrophenol	μg/L	-	-	3	0	<10	-	19	0	<100	-					
121-14-2	2,4-Dinitrotoluene	μg/L	-	-	9	0	<20	-	51	0	<20	-					
606-20-2	2,6-Dinitrotoluene	μg/L	-	-	9	0	<20	-	51	0	<20	-					
35572-78-2	2-Amino-4,6-Dinitrotoluene	μg/L	-	-	3	0	<5	-	13	0	<5	=					
78-93-3	2-Butanone	μg/L	-	-	4	0	<5	-	19	0	<5	-					
110-75-8	2-Chloroethyl Vinyl Ether	μg/L	-	-	5	0	<1	-	19	0	<1	-					
91-58-7	2-Chloronaphthalene	μg/L	-	-	3	0	<1	-	19	0	<10	-					
95-57-8	2-Chlorophenol	μg/L	-	-	3	0	<1	-	19	0	<10	-					
95-49-8	2-Chlorotoluene	μg/L	-	140	34	0	<0.5	-	475	0	<1	-			/		
591-78-6	2-Hexanone	μg/L	-	-	4	0	<5	-	19	0	<5	-					
91-57-6	2-Methylnaphthalene	μg/L	-	-	3	0	<1	-	19	0	<10	-					
95-48-7	2-Methylphenol	μg/L	-	-	3	0	<1	-	19	0	<10	-					
88-74-4	2-Nitroaniline	μg/L	-	-	3	0	<1	-	19	0	<10	-					
88-75-5	2-Nitrophenol	μg/L	_	-	3	0	<1	-	19	0	<10	_					
88-72-2	2-Nitrotoluene	μg/L	-	-	3	0	<5	_	13	0	<5	_					
7642-04-8	2-Octene, (Z)-	μg/L	_	_	1	1	34	_	5	5	250	_	<b>√</b>	<b>√</b>			
84989-04-8	3 & 4-Methylphenol	μg/L	_	_	3	0	<1	_	19	0	<10	_					
91-94-1	3,3'-Dichlorobenzidine	μg/L	_	_	3	0	<5	_	19	0	<50	_				1	
95-76-1	3,4-Dichloroaniline	μg/L	_	_	3	0	<1	_	13	0	<1	_				1	
51-36-5	3,5-Dichlorobenzoic acid	μg/L	_	_	3	0	<1	_	19	0	<1	_				1	
16655-82-6	3-Hydroxycarbofuran	μg/L	_	_	5	0	<10	_	19	0	<2	_				_	
99-09-2	3-Nitroaniline	μg/L μg/L	_	_	3	0	<1	_	19	0	<10	_				_	
99-08-1	3-Nitrotoluene	μg/L μg/L	_	_	3	0	<5	_	13	0	<5	_					
72-54-8	4,4´-DDD	μg/L μg/L		_	8	0	<1		38	0	<1					+	
72-55-9	4,4´-DDE	μg/L μg/L		_	8	0	<2	_	38	0	<2	-				+	
50-29-3	4,4´-DDT	μg/L μg/L	_	_	8	0	<1		38	0	<1					+	4
534-52-1	4,6-Dinitro-2-Methyl phenol	μg/L μg/L		_	3	0	<5		19	0	<50					+	
19406-51-0	4-Amino-2,6-Dinitrotoluene		-	_	3	0	<5	_	13	0	<5					+	+
101-55-3	4-Ammo-2,0-Dintrotoidene 4-Bromophenyl Phenyl Ether	μg/L	-	_	3	0	<1	-	19	0	<10					-	4
59-50-7	4-Chloro-3-Methylphenol	μg/L		-	3	0	<1	-	19	0	<10	-				_	4
	, ,	μg/L	-	-	3	_		-		-		-				-	
106-47-8	4-Chloroaniline	μg/L	-	-	3	0	<1	-	19	0	<10	-				-	
7005-72-3	4-Chlorophenyl Phenyl Ether	μg/L		140	3	0	<1	-	19	0	<10	-			/	+	4
106-43-4	4-Chlorotoluene	μg/L	-	140	34	0	<0.5	-	475	0	<1	-			<i>/</i>	+	4
108-10-1	4-Methyl-2-pentanone	μg/L	-	120	34	0	<5	-	19	0	<5	-			v		4—
100-01-6	4-Nitroaniline	μg/L	-	-	3	0	<1	-	19	0	<10	-					1
100-02-7	4-Nitrophenol	μg/L	-	-	3	0	<5	-	19	0	<50	-					4—
99-99-0	4-Nitrotoluene	μg/L	-	-	3	0	<5	-	13	0	<5	-					4
104-40-5	4-Nonylphenol	μg/L	-	-	3	0	<0.025	-	13	0	<0.25	-					4
1806-26-4	4-Octylphenol	μg/L	-	-	3	0	<0.025	-	13	0	<0.25	-					4
140-66-9	4-tert-Octylphenol	μg/L	-	-	3	0	<0.005	-	13	1	0.014	-					4
2315-61-9	4-tert-Octylphenol diethoxylate	μg/L	-	-	3	0	<0.025	-	13	0	<0.25	-					
1173019-48-1	4-tert-Octylphenol monoethoxylate	μg/L	-	-	3	0	<0.025	-	13	4	0.072	-		✓			
756426-58-1	9CI-PF3ONS	μg/L	-	-	8	2	0.0017	-	16	0	<0.002	-	✓				
83-32-9	Acenaphthene	μg/L	-	-	6	0	<5	-	38	0	<10	-					
208-96-8	Acenaphthylene	μg/L	-	-	6	0	<5	-	38	0	<10	-					







						Producti	on Wells			Monitor	ing Wells			Meets	COPC Crite	Screen ia*	ing
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Initio	il 3 4	Seco 5	ondary 6
75-07-0	Acetaldehyde	μg/L	-	-	3	0	<2	-	14	2	0.48	-					
103-90-2	Acetaminophen	μg/L	-	-	3	0	<0.18	-	13	0	<0.32	-					
34256-82-1	Acetochlor	μg/L	-	-	3	0	<1	-	19	0	<1	-					
187022-11-3	Acetochlor ESA	μg/L	-	-	3	0	<0.03	-	13	0	< 0.03	-					
194992-44-4	Acetochlor OA	μg/L	-	-	3	1	0.0068	-	13	0	< 0.033	-	✓				
67-64-1	Acetone	μg/L	-	-	4	2	790	-	16	10	31	-	✓	✓			
75-05-8	Acetonitrile	μg/L	-	-	4	2	7.9	-	19	7	0.62	-	✓	✓			
50594-66-6	Acifluorfen	μg/L	-	-	5	0	<0.4	-	19	0	<0.4	-					
107-02-8	Acrolein	μg/L	-	-	4	0	<5	-	20	0	<5	-					
107-13-1	Acrylonitrile	μg/L	-	-	4	0	<2	-	20	0	<2	-					
958445-44-8	ADONA	μg/L	-	-	8	2	0.0017	-	16	0	<0.002	-	✓				
15972-60-8	Alachlor	μg/L	2	-	15	0	<1	-	19	0	<1	-			/		
142363-53-9	Alachlor ESA	μg/L	-	-	3	0	<0.042	-	13	1	0.021	-					
171262-17-2	Alachlor OA	μg/L	-	-	3	0	<0.033	-	13	0	< 0.033	-					
116-06-3	Aldicarb	μg/L	-	-	5	0	<10	-	19	0	<2	-					
1646-88-4	Aldicarb Sulfone	μg/L	-	-	5	0	<10	-	19	0	<2	-					
1646-87-3	Aldicarb sulfoxide	μg/L	-	-	5	0	<10	-	19	0	<2	-					
309-00-2	Aldrin	μg/L	-	-	12	0	<1	-	38	0	<1	-					
ALK	Alkalinity (as CaCO3)	μg/L	-	-	6	6	430000	-	35	35	430000	-	✓	✓			
ALKC-1	Alkalinity, Carbonate as CaCO3	μg/L	-	-	15	6	401000	-	19	2	63000	-	✓				
ALKH-1	Alkalinity, Hydroxide as CaCO3	μg/L	-	-	9	0	<5000	-	19	0	<5000	-					
107-05-1	Allyl Chloride	μg/L	-	-	4	0	<0.5	-	20	0	<0.5	-					
319-84-6	alpha-BHC	μg/L	-	-	8	0	<1	-	38	0	<1	-					
5103-71-9	alpha-Chlordane	μg/L	-	-	3	0	<1	-	19	0	<1	-					
7429-90-5	Aluminum	μg/L	1000	-	12	7	2400	2.40	38	38	2800	2.80	✓	✓	/ /	✓	
7664-41-7	Ammonia	μg/L	-	-	3	1	34	-	13	4	1200	-	✓	✓			
26787-78-0	Amoxicillin	μg/L	-	-	3	1	0.0055	-	13	0	< 0.01	-	✓				
62-53-3	Aniline	μg/L	-	-	3	0	<1	-	19	0	<10	-					
120-12-7	Anthracene	μg/L	-	-	6	0	<5	-	38	0	<10	-					
7440-36-0	Antimony	μg/L	6	-	12	6	0.2	0.03	38	37	0.39	0.07	✓	✓	✓ ✓		
12674-11-2	Aroclor 1016	μg/L	-	-	3	0	<0.1	-	19	0	<25	-					
11104-28-2	Aroclor 1221	μg/L	-	-	3	0	<0.1	-	19	0	<25	-					
11141-16-5	Aroclor 1232	μg/L	-	-	3	0	<0.1	-	19	0	<25	-					
53469-21-9	Aroclor 1242	μg/L	-	-	3	0	<0.1	-	19	0	<25	-					
12672-29-6	Aroclor 1248	μg/L	-	-	3	0	<0.1	-	19	0	<25	-					
11097-69-1	Aroclor 1254	μg/L	-	-	3	0	<0.1	-	19	0	<25	-					
11096-82-5	Aroclor 1260	μg/L	-	-	3	0	<0.1	-	19	0	<25	-					
7440-38-2	Arsenic	μg/L	10	-	12	11	2	0.20	41	39	38	3.80	✓	<b>√</b>	/ /	✓	
22541-54-4	Arsenic III	μg/L	-	-	6	2	0.17	-	38	14	5.4	-	✓	✓			
17428-41-0	Arsenic V	μg/L	-		6	6	1.4	-	38	36	1.5	-	✓	✓			
1332-21-4	Asbestos	MFL	7	-	3	0	<1.7	-	19	0	<5	-			/		
29122-68-7	Atenolol	μg/L	-	-	3	1	0.00028	-	13	2	0.0066	-	✓				
134523-00-5	Atorvastatin	μg/L	-	-	3	1	0.0011	-	13	1	0.00076	-	✓				
1912-24-9	Atrazine	μg/L	1	-	15	0	<1	-	19	0	<1	-			/		
83905-01-5	Azithromycin	μg/L	-	-	3	1	0.0057	-	13	2	0.015	-	$\checkmark$				







According   Parameter   Units   MCL   NL, Number of Disservation   Disservation							Producti	on Wells			Monitor	ing Wells			Meet	COPC	: Screen ria*	ing
7440-39-31   Вашит   руд. 1000   12   12   78   0.00   38   38   100   0.10   7   7   7	CASRN	Parameter	Units	MCL	NL			Observed	Maximum to			Observed	Maximum to	1				ndary 6
2007-09-09   Bentacon	103-33-3	Azobenzene	μg/L	-	-	2	0	<1	-	15	0	<10	-					
100-52-7	7440-39-3	Barium	μg/L	1000	-	12	12	76	0.08	38	38	100	0.10	✓	✓	✓ ✓		
Principle   Prin	25057-89-0	Bentazon	μg/L	18	-	15	0	<2	-	19	0	<2	-			✓		
\$2.75	100-52-7	Benzaldehyde	μg/L	-	-	3	0	<2	-	14	0	<2	-					
September   Benzolamintacene   Ug/L   -   -   6   0   -   3   -   38   0   -   10   -	71-43-2	Benzene	μg/L	1	-	201	0	<0.5	-	475	24	19	19.00		✓	✓ ✓	✓	✓
Solution   Solution	92-87-5	Benzidine	μg/L	-	-	3	0	<10	-	19	0	<100	-					
2015-59-2   Berosphthoconthene   1914	56-55-3	Benzo[a]anthracene	μg/L	-	-	6	0	<5	-	38	0	<10	-					
1973-42   Beronoj(a))perpriene	50-32-8	Benzo[a]pyrene	μg/L	0.2	-	14	0	<1	-	38	0	<10	-			✓		
2017-08-9	205-99-2	Benzo[b]fluoranthene	μg/L	-	-	6	0	<5	-	38	0	<10	=					
6:85-0   Berroti Rold   Ug/L   3   0   <100   -   19   2   6.7	191-24-2	Benzo[g,h,i]perylene	μg/L	-	-	6	0	<5	-	38	0	<20	=					
Section   Sect	207-08-9	Benzo[k]fluoranthene	μg/L	-	-	6	0	<5	-	38	0	<10	-					
100-51-6	65-85-0	Benzoic Acid	μg/L	-	-	3	0	<100	-	19	2	6.7	-					
Bensyl burly phthalate	95-16-9	Benzothiazole	μg/L	-	-	0	0	-	-	2	2	8.5	-					
TH4O-41-7   Beryllium	100-51-6	Benzyl Alcohol	μg/L	-	-	3	0	<1	-	19	0	<10	-					
319-85-7     bets-BHC   190/L   -       -       -     -     -     -	85-68-7	Benzyl butyl phthalate	μg/L	-	-	6	2	3.7	-	38	10	6.3	-	✓	✓			
319-85-7	7440-41-7	Beryllium	μq/L	4	-	12	1	0.088	0.02	38	1	0.095	0.02	✓		✓		
Principal	319-85-7	beta-BHC		-	-	8	0	<2	-	38	0	<2						
BOD   Biochemical Oxygen Demand   μg/L     3   0   <2000   .   13   2   6900   .	71-52-3	Bicarbonate Alkalinity as HCO3		-	-	9	9	520000	-	19	19	520000	-	✓	✓			
111-91-1   bis(2-chloreothoxy)methane   µg/L   -   3   0   x1   -   19   0   x10   -	BOD			-	-	3	0	<2000	-	13	2	6900	-					
111-44-4   bis(2-chlorospropybether   µg/L   -   3   0   41   -   19   0   410   -	111-91-1			-	-	3	0	<1	-		0		-					
103-60-1   bis(2-chtoriospropy))ether   µg/L   -				_	-	3	0	<1	-		0		-					
103-23-1   Bis(2-ethylhevyl)aclipate   µg/L   400   -   11   0   <50   -   19   0   <50   -	108-60-1			_	-	3	0	<1	-		0		-					
117-81-7				400	_	11	0	<50	-		0		-			<b>✓</b>		
80-05-7   Bisphenol A					_	18			-		5		0.23		✓	<b>✓ ✓</b>		
7440-42-8         Boron         µg/L         -         1000         7         7         200         0.20         43         41         590         0.59         ✓         <				-	_		3		-		13		-	✓	✓			
314-40-9   Bromacil		'		-	1000	7	7		0.20				0.59	<b>√</b>	✓	/ /	<b>✓</b>	
15541-45-4   Bromate				_		5	0		_									
24959-67-9         Bromide         μg/L         -         -         3         3         850         -         14         14         5700         -         ✓         ✓           108-86-1         Bromocherene         μg/L         -         -         34         0         <0.5		1		10	_	3	_		-		6		16.00		✓	/ /	<b>✓</b>	
108-86-1   Bromobenzene				-	_	3	_		_					<b>√</b>	<b>√</b>			
S589-96-8   Bromochloroacetic acid   μg/L   -   -   3   0   <1   -   19   0   <1   -				_	_	34			_				_					
74-97-5   Bromochloromethane   μg/L   -   -   34   0   <0.5   -   475   0   <1   -				-	-	3	_		_		-	_	_					
71133-14-7   Bromodichloroacetic acid   μg/L   -   -   0   0   -   -   6   0   <1   -						34	_											
75-27-4         Bromodichloromethane         µg/L         -         201         3         0.5         -         475         4         1.5         -         ✓         ✓         Image: Contract of the co													1					
75-25-2   Bromoform   μg/L   -   -   196   2   1.5   -   475   0   <1   -						-			_					<b>✓</b>	<b>√</b>			
74-83-9         Bromomethane         μg/L         -         -         34         0         <0.5         -         475         0         <1         -         Image: square squar									_		•							
23184-66-9         Butachlor         μg/L         -         -         5         0         <1         -         19         0         <1         -         -         Image: stress of the content of the co									_									
123-72-8     Butanal     μg/L     -     -     3     0     <2																		
7440-43-9         Cadmium         μg/L         5         -         12         5         0.23         0.05         38         16         0.12         0.02         ✓									_									
58-08-2         Caffeine         μg/L         -         -         6         3         0.031         -         32         15         0.13         -         ✓         ✓           7440-70-2         Calcium         μg/L         -         -         12         12         183000         -         32         32         178000         -         ✓         ✓           133-06-2         Captan         μg/L         -         -         5         0         <10					_		_		0.05					<b>✓</b>	<b>√</b>	<b>✓</b>		
7440-70-2     Calcium     μg/L     -     -     12     12     183000     -     32     32     178000     -     √     √       133-06-2     Captan     μg/L     -     -     5     0     <10				-	_				-									
$\begin{array}{c ccccccccccccccccccccccccccccccccccc$				_	_				_									
298-46-4     Carbamazepine     μg/L     -     -     3     1     0.00009     -     13     0     <0.001     -     ✓     -       63-25-2     Carbaryl     μg/L     -     -     5     0     <10																		
63-25-2 Carbaryl µg/L 5 0 <10 - 19 0 <2 -						- J								<b>/</b>				1
							'						+					
86-74-8   Carbazole   ug/    -   -   3   0   <1   -   19   0   <10   -	86-74-8	Carbaryi	μg/L μg/L			3	0	<10		19	0	<10						4







						Producti	on Wells		Monitoring Wells						Meets COPC Screening Criteria*					
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Init 2	ial 3		Secondar 5 6	<i>y</i>		
1563-66-2	Carbofuran	μg/L	18	-	15	0	<10	-	19	0	<2	-			✓					
75-15-0	Carbon Disulfide	μg/L	-	160	32	0	<0.5	-	19	1	0.26	0.00			✓					
56-23-5	Carbon Tetrachloride	μg/L	0.5	-	201	2	0.15	0.30	475	8	1	2.00	✓	✓	✓	✓	✓ ✓	/		
133-90-4	Chloramben	μg/L	-	-	3	0	<1	-	19	0	<1	-								
7775-09-9	Chlorate	μg/L	-	-	3	2	93	-	13	12	160	-	✓	✓						
57-74-9	Chlordane	μg/L	0.1	-	15	0	<0.1	-	19	0	<5	-			✓					
16887-00-6	Chloride	μg/L	-	-	9	9	130000	-	24	24	601000	-	<b>✓</b>	✓						
76-13-1	Chlorinated Fluorocarbon (Freon 113)	μg/L	1200	-	34	0	<10	-	19	0	<5	-			✓					
7782-50-5RF	Chlorine Residual, Free	μg/L	-	-	3	3	35	-	13	10	80	-	✓	✓						
TOTAL-CHLORINE	Chlorine Residual, Total	μg/L	-	-	3	3	48	-	13	13	160	-	✓	✓						
14998-27-7	Chlorite	μg/L	1000	-	3	2	14	0.01	19	7	16	0.02	✓	✓	✓					
79-11-8	Chloroacetic Acid	μg/L	-	-	3	0	<2	-	19	0	<2	-		1	i					
108-90-7	Chlorobenzene	μg/L	70	-	34	0	<0.5	-	475	0	<1	-		1	✓					
5278-95-5	Chlorodibromoacetic acid	μg/L	-	-	0	0	-	-	6	0	<1	-			i					
75-00-3	Chloroethane	μg/L	-	-	34	0	<0.5	-	475	0	<1	-			i					
67-66-3	Chloroform	μg/L	-	-	201	188	5.4	-	475	182	37	-	✓	✓	i					
74-87-3	Chloromethane	μg/L	-	-	34	0	<0.5	-	475	0	<1	-			i					
1897-45-6	Chlorothalonil	μg/L	_	-	5	0	<5	-	19	0	<2.5	-								
101-21-3	Chlorpropham	μg/L	_	-	5	0	<1	-	19	0	<1	-								
7440-47-3	Chromium	μg/L	50	-	13	8	7.2	0.14	38	31	290	5.80	✓	<b>√</b>	✓	✓	✓			
18540-29-9	Chromium, Hexavalent	μg/L	_	-	5	5	1.9	-	19	13	5.9	-	✓	<b>√</b>						
218-01-9	Chrysene	μg/L	_	-	6	0	<5	-	38	0	<10	-								
85721-33-1	Ciprofloxacin	μg/L	_	-	3	2	0.11	-	13	11	0.014	-	✓	<b>√</b>						
156-59-2	cis-1,2-Dichloroethene	μg/L	6	-	201	110	3.4	0.57	475	93	300	50.00	✓	<b>√</b>	✓	✓	<b>√</b> ✓	/		
10061-01-5	cis-1,3-Dichloropropene	μg/L	_	-	34	0	<0.5	-	475	0	<1	-								
5103-73-1	cis-Nonachlor	μg/L	_	-	3	0	<0.1	-	18	0	<0.1	-								
7440-48-4	Cobalt	μg/L	_	_	6	6	1	-	26	25	8.3	-	<b>√</b>	<b>√</b>	i					
CLR	Color	Units	_	-	9	6	5	-	19	0	<3	-	✓							
7440-50-8	Copper	μg/L	1300	_	12	8	49	0.04	38	29	33	0.03	<b>√</b>	<b>√</b>	<b>√</b>			$\overline{}$		
486-56-6	Cotinine	μg/L	_	-	3	1	0.0012	-	13	11	0.027	-	<b>√</b>	<b>√</b>	i			-		
123-73-9	Crotonaldehyde	μg/L	_	_	3	0	<2	-	14	0	<2	-		,	i					
21725-46-2	Cyanazine	<u>μg</u> /L	_	_	5	0	<1	-	19	0	<1	-		,	i			-		
57-12-5	Cyanide	μg/L	150	-	9	0	<100	-	19	0	<5	-		,	✓					
1122-82-3	Cyclohexane, isothiocyanato-	μg/L	-	_	0	0	-	-	1	1	1	-		,	i					
108-94-1	Cyclohexanone	<u>μg/L</u>	_	_	3	0	<2	-	14	0	<2	_		, 1	1			-		
75-99-0	Dalapon	<u>μg</u> /L	200	_	11	0	<10	-	19	0	<0.4	-		1	<b>√</b>					
1861-32-1	DCPA	<u>μ</u> g/L	-	_	5	0	<0.1	-	19	0	<0.1	_		,				-		
112-31-2	Decanal	<u>μ</u> g/L	-	_	3	0	<2	-	14	0	<2	-		1				$\overline{}$		
134-62-3	DEET	μg/L	_	_	3	3	0.0043	_	13	13	0.0036	_	<b>✓</b>	<b>√</b>						
319-86-8	delta-BHC	μg/L μg/L	_	_	8	0	<1	-	38	0	<1	_		,						
439-14-5	Diazepam	μg/L μg/L	_	_	3	0	<0.001	_	13	1	0.00021	_		,				<b>—</b>		
333-41-5	Diazinon	μg/L μg/L	_	1.2	5	0	<1	-	19	1	0.055	0.05		$\rightarrow$	<b>√</b>					
53-70-3	Dibenzo[a,h]anthracene	μg/L μg/L	_	-	6	0	<5	-	38	0	<20	-		,						
132-64-9	Dibenzofuran	μg/L μg/L	_	_	3	0	<1	-	19	0	<10	_		,						
631-64-1	Dibromoacetic acid	μg/L μg/L	_	_	3	0	<1	-	19	0	<1	_		,						







						Producti	on Wells			Monitor	ing Wells		Meets COPC Scree Criteria*				ing
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Initia 2	al 4	Seco	ondary 6
124-48-1	Dibromochloromethane	μg/L	-	-	201	1	0.6	-	475	1	70	-	✓				
74-95-3	Dibromomethane	μg/L	-	-	34	0	<0.5	-	475	0	<1	-					
1918-00-9	Dicamba	μg/L	-	-	5	0	<1.5	-	19	0	<0.6	-					
3400-09-7	Dichloramine	μg/L	-	-	3	3	7	-	13	10	15	-	✓	✓			
79-43-6	Dichloroacetic Acid	μg/L	-	-	3	0	<1	-	19	0	<1	-					
75-71-8	Dichlorodifluoromethane	μg/L	-	1000	34	0	<0.5	-	475	0	<1	=		`			
120-36-5	Dichloroprop	μg/L	-	-	5	0	<0.3	-	19	0	<0.3	-					
15307-86-5	Diclofenac	μg/L	-	-	3	0	<0.001	-	13	2	0.00073	=					
60-57-1	Dieldrin	μg/L	-	-	12	0	<2	-	38	0	<2	=					
84-66-2	Diethyl phthalate	μg/L	-	-	6	0	<20	-	38	16	0.88	-		✓			
56-53-1	Diethylstilbestrol	μg/L	-	-	3	0	<0.005	-	13	0	<0.1	-					
108-20-3	Di-isopropyl ether	μg/L	-	-	172	0	<3	-	473	13	6.4	-		✓			
60-51-5	Dimethoate	μg/L	-	-	5	0	<2	-	19	0	<2	-					
131-11-3	Dimethyl phthalate	μg/L	-	-	6	0	<20	-	38	0	<20	-					
84-74-2	Di-n-butyl-phthalate	μg/L	-	-	6	2	0.23	-	38	10	2.7	-	✓	✓			
117-84-0	Di-n-octyl phthalate	μg/L	-	-	6	0	<5	-	38	0	<10	-					
88-85-7	Dinoseb	μg/L	7	-	15	0	<2	-	19	0	<0.4	-		`			
957-51-7	Diphenamid	μg/L	-	-	5	0	<100	-	19	0	<1	-					
85-00-7	Diquat	μg/L	20	-	15	0	<13	-	19	0	<32	-		`	/		
298-04-4	Disulfoton	μg/L	-	-	5	0	<1	-	19	0	<2	-					
330-54-1	Diuron	μg/L	-	-	3	0	<1	-	13	0	<1	-					
959-98-8	Endosulfan I	μg/L	-	-	8	0	<10	-	38	0	<10	-					
33213-65-9	Endosulfan II	μg/L	-	-	8	0	<2	-	38	0	<2	=					
1031-07-8	Endosulfan sulfate	μg/L	-	-	8	0	<2	-	38	0	<2	=					
145-73-3	Endothall	μg/L	100	-	15	0	<45	-	19	0	<45	-		`		j	
72-20-8	Endrin	μg/L	2	-	18	0	<2	-	38	0	<2	-		`			
7421-93-4	Endrin Aldehyde	μg/L	-	-	8	0	<2	-	38	0	<2	-					
53494-70-5	Endrin Ketone	μg/L	-	-	3	0	<1	-	19	0	<1	-					
481-30-1	Epitestosterone	μg/L	-	-	3	0	<0.001	-	13	0	< 0.001	-					
759-94-4	EPTC	μg/L	-	-	5	0	<1	-	19	0	<1	-					
50-27-1	Estriol	μg/L	-	-	3	0	<0.005	-	13	0	<0.16	-					
53-16-7	Estrone	μg/L	-	-	3	0	<0.001	-	13	3	0.00077	-		✓			
2991-50-6	EtFOSAA	μg/L	-	-	6	0	<0.002	_	16	1	0.0016	-					
64-17-5	Ethanol	μg/L	-	-	3	0	<10000	_	13	0	<10000	-					
563-12-2	Ethion	μg/L	-	-	3	0	<1	-	19	0	<1	-					
97-63-2	Ethyl Methacrylate	μg/L	-	-	4	0	<5	-	20	0	<5	-					
637-92-3	Ethyl tert-butyl ether	μg/L	-	-	201	0	<3	-	473	0	<2	-					
100-41-4	Ethylbenzene	μg/L	300	-	201	0	<0.5	-	475	0	<1	-		,	/		
107-21-1	Ethylene Glycol	μg/L	-	14000	3	0	<10000	-	13	0	<10000	-		,	/		
206-44-0	Fluoranthene	μg/L	-	-	6	0	<5	-	38	0	<10	-					
86-73-7	Fluorene	μg/L	-	-	6	0	<5	-	38	0	<10	-					
16984-48-8	Fluoride	μg/L	2000	-	9	9	390	0.20	19	19	370	0.19	✓	√ \	/ /		
54910-89-3	Fluoxetine	μg/L	-	_	3	2	0.0011	-	13	8	0.0011	-	✓	✓			
944-22-9	Fonofos	μg/L	-	-	3	0	<0.015	-	13	0	< 0.035	-					
50-00-0	Formaldehyde	μg/L	-	100	3	0	<2	-	14	1	0.89	0.01		`			







					Production Wells Monitoring Wells									Meets COPC Screen Criteria* Initial Seco			
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Init 2	ial 3		Secondary 5 6
1222-05-5	Galaxolide (HHCB)	μg/L	-	-	3	3	0.16	-	13	12	0.057	-	✓	✓			
58-89-9	gamma-BHC	μg/L	0.2	-	18	0	<1	-	38	0	<1	-			✓		
5566-34-7	gamma-Chlordane	μg/L	-	-	3	0	<1	-	19	0	<1	-					
8032-32-4	Gasoline Range Organics	μg/L	-	-	3	2	160	=	13	11	860	-	✓	✓			
25812-30-0	Gemfibrozil	μg/L	-	-	3	1	0.000098	-	13	2	0.00012	-	✓				
107-22-2	Glyoxal	μg/L	-	-	3	0	<2	=	14	0	<2	-					
1071-83-6	Glyphosate	μg/L	700	-	15	0	<25	-	19	2	2.2	0.00			<b>✓</b>		
12587-46-1	Gross Alpha	pci/L	15	-	7	6	11	0.73	19	19	7.94	0.53	✓	✓	✓	✓ ,	✓
12587-47-2	Gross Beta	pci/L	-	-	3	3	8.1	-	19	19	12	-	✓	✓			
HAA5	HAA5, Total	μg/L	60	-	3	0	<1	-	19	0	<1	-			✓		
Hardt-1	Hardness, total as CaCO3	μg/L	-	-	12	12	834000	-	30	30	775000	-	✓	✓			
76-44-8	Heptachlor	μg/L	0.01	-	12	0	<1	-	38	0	<1	-			✓		
1024-57-3	Heptachlor Epoxide	μg/L	0.01	-	14	0	<1	-	38	0	<1	-			✓		
111-71-7	Heptanal	μg/L	-	-	3	0	<2	-	14	0	<2	-					
118-74-1	Hexachlorobenzene	μg/L	1	-	17	0	<1	-	57	3	0.09	0.09		✓	✓	✓	
87-68-3	Hexachlorobutadiene	μq/L	-	-	37	0	<1	=	494	0	<10	-					
77-47-4	Hexachlorocyclopentadiene	μg/L	50	-	17	0	<10	=	57	0	<50	-			✓		
67-72-1	Hexachloroethane	μg/L	-	_	3	0	<1	-	19	0	<10	-					
66-25-1	Hexanal	μg/L	_	_	3	0	<2	-	14	0	<2	-					
13252-13-6	HFPO-DA	μg/L	-	_	6	0	<0.002	-	16	0	<0.002	-					
2691-41-0	HMX	μg/L	-	350	3	0	<5	=	13	0	<5	-			✓		
302-01-2	Hydrazine	μg/L	-	-	3	0	<1	=	13	0	<1	-					
15687-27-1	Ibuprofen	μg/L	-	_	3	1	0.0013	-	13	9	0.0068	-	✓	✓			
193-39-5	Indeno[1,2,3-cd]pyrene	µg/L	-	_	6	0	<5	-	38	0	<20	-					
20461-54-5	lodide	μg/L	-	_	3	3	34	-	13	11	76	-	<b>✓</b>	<b>√</b>			
74-88-4	Iodomethane	μg/L	-	-	4	0	<0.5	=	20	0	<0.5	-					
73334-07-3	lopromide	μg/L	-	_	3	0	<0.006	-	13	0	<0.08	-					
7439-89-6	Iron	μg/L	-	-	13	9	3200	-	35	26	6600	-	✓	✓			
78-59-1	Isophorone	μg/L	-	_	3	0	<1	-	19	0	<10	-					
67-63-0	Isopropyl Alcohol	μg/L	-	_	3	0	<10000	-	13	0	<10000	-					
98-82-8	Isopropylbenzene	μg/L	_	770	34	0	<0.5	-	475	0	<1	-			<b>✓</b>		
7439-92-1	Lead	μg/L	15	-	12	6	4	0.27	32	14	1.2	0.08	✓	✓	<b>√</b>	<b>√</b>	
330-55-2	Linuron	μg/L	-	_	3	0	<1	-	13	0	<1	-					
7439-93-2	Lithium	μg/L	_	_	6	0	<10	-	26	2	25	-					
179601-23-1	m,p-Xylene	μg/L	-	_	201	0	<0.5	-	19	0	<0.5	-					
7439-95-4	Magnesium	μg/L	-	_	12	12	91500	-	32	32	102000	-	<b>✓</b>	<b>√</b>			
7439-96-5	Manganese	μg/L	-	500	13	13	67	0.13	38	38	340	0.68	✓	✓	<b>√</b>	✓ ,	✓
2355-31-9	MeFOSAA	μg/L	-	-	6	0	<0.002	-	16	1	0.0013	-					
57-53-4	Meprobamate	μg/L	_	_	3	0	<0.001	-	13	<u>.</u> 1	0.0022	-					
7439-97-6	Mercury	μg/L	2	_	12	4	0.023	0.01	38	12	0.025	0.01	<b>√</b>	<b>√</b>	<b>√</b>		
126-98-7	Methacrylonitrile	μg/L	-	_	4	0	<0.5	-	20	0	<0.5	-					
76-99-3	Methadone	μg/L	_	-	3	1	0.00016	-	13	5	0.00024	_	<b>✓</b>	<b>√</b>			
67-56-1	Methanol	μg/L	_	-	3	0	<10000	-	13	0	<10000	_					
2032-65-7	Methiocarb	μg/L	_	_	5	0	<10	-	19	0	<2	_					
16752-77-5	Methomyl	μg/L	_	_	5	0	<10	_	19	0	<2	_					







			MC			Producti	on Wells			Monitor	ing Wells			Meets	COPC Crite	Screen	ing
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Initio	il 3 4	Seco 5	ondary 6
72-43-5	Methoxychlor	μg/L	30	-	18	0	<10	-	38	0	<2	-		,	/		
78-98-8	Methyl Glyoxal	μg/L	-	-	3	0	<2	_	14	0	<2	-					
80-62-6	Methyl Methacrylate	μg/L	-	_	4	0	<0.5	_	20	2	0.13	-					
1634-04-4	Methyl tert-butyl ether (MTBE)	μg/L	13	-	200	0	<3	_	475	18	67	5.15		✓ ,	/ /	✓	✓
MBAS	Methylene blue active substances (MBAS)	μg/L	-	-	9	5	100	-	19	12	83	-	<b>✓</b>	<b>√</b>			
75-09-2	Methylene Chloride	μg/L	5	-	201	0	<0.5	-	475	6	2	0.40		✓ ,	/ /		
51218-45-2	Metolachlor	μg/L	-	-	5	0	<1	-	19	0	<1	-					
171118-09-5	Metolachlor ESA	μg/L	-	-	3	0	<0.01	-	13	0	< 0.01	-					
152019-73-3	Metolachlor OA	μg/L	-	-	3	1	0.0039	-	13	0	< 0.01	-	✓				
21087-64-9	Metribuzin	μg/L	-	-	5	0	<1	-	19	1	0.079	-				j	
2385-85-5	Mirex	μg/L	-	-	3	0	<0.1	-	18	0	<0.1	-					
2212-67-1	Molinate	μg/L	20	-	11	0	<2	-	19	0	<1	=		,			
7439-98-7	Molybdenum	μg/L	-	-	10	10	9.8	-	26	26	60	-	✓	✓			
79-08-3	Monobromoacetic acid	μg/L	-	-	3	0	<1	-	19	0	<1	-					
10599-90-3	Monochloramine	μg/L	-	-	3	3	6	-	13	13	100	-	✓	✓			
60-34-4	Monomethylhydrazine	μg/L	-	-	3	0	<2	-	13	0	<2	-				j	
57-27-2	Morphine	μg/L	-	-	3	1	0.00058	-	13	10	0.0018	-	✓	✓			
91-20-3	Naphthalene	μg/L	-	17	40	0	<5	-	513	0	<10	-		,			
22204-53-1	Naproxen	μg/L	-	-	3	1	0.0011	-	13	2	0.0029	=	✓				
104-51-8	n-Butylbenzene	μg/L	-	260	34	0	<0.5	-	475	0	<1	-		,			
7440-02-0	Nickel	μg/L	100	-	12	6	4.1	0.04	38	38	200	2.00	✓	✓ ,	/ /	✓	
14797-55-8	Nitrate as N	μg/L	10000	-	45	45	7400	0.74	21	17	16000	1.60	✓	✓ ,	/ /	✓	
NN	Nitrate plus nitrite as N	μg/L	10000	-	7	7	5800	0.58	21	17	16000	1.60	✓	✓ ,	/ /	✓	
14797-65-0	Nitrite as N	μg/L	1000	-	9	0	<100	-	21	10	160	0.16		✓ ,	/ /		
98-95-3	Nitrobenzene	μg/L	-	-	10	0	<10	-	52	0	<10	-				j	
55-18-5	n-Nitrosodiethylamine	μg/L	-	0.01	3	1	0.0015	0.15	19	1	0.0011	0.11	✓	,	/ /		
62-75-9	N-Nitrosodimethylamine	μg/L	-	0.01	6	0	<1	-	38	1	0.0018	0.18		,			
924-16-3	n-Nitrosodi-n-butylamine	μg/L	-	-	3	0	<0.002	-	19	0	<0.002	-					
621-64-7	n-Nitrosodi-n-propylamine	μg/L	-	0.01	6	0	<1	-	38	0	<10	-		,			
86-30-6	n-Nitrosodiphenylamine	μg/L	-	-	3	0	<1	-	19	0	<10	-					
10595-95-6	N-Nitrosomethylethylamine	μg/L	-	_	3	0	<0.002	-	19	0	< 0.002	-					
59-89-2	n-Nitrosomorpholine	μg/L	-	-	3	2	0.039	-	19	11	0.035	-	✓	✓			
100-75-4	n-Nitrosopiperidine	μg/L	-	-	3	0	<0.002	-	19	0	<0.002	-					
930-55-2	n-Nitrosopyrrolidine	μg/L	-	-	3	0	<0.002	-	19	0	<0.002	-					
25154-52-3	Nonylphenol	μg/L	-	-	3	1	0.016	-	13	4	4.4	-	✓	✓			
20427-84-3	Nonylphenol diethoxylate	μg/L	-	-	3	1	0.05	-	13	3	3.2	-	✓	✓			
27986-36-3	Nonylphenol monoethoxylate	μg/L	-	-	3	1	0.011	-	13	9	2.3	-	✓	✓			
103-65-1	n-Propylbenzene	μg/L	-	260	34	0	<0.5	-	475	0	<1	-		,	/		
136777-61-2	o&p-Xylene	μg/L	-	-	0	0	-	-	456	0	<1	-					
111-66-0	Octene-1	μg/L	-	-	0	0	-	-	1	1	25	-					
OILGREASE	Oil & Grease (HEM)	μg/L		-	3	2	3100	-	13	2	1500	-	✓				
14265-44-2	o-Phosphate as P	μg/L	-	-	3	3	150	-	13	13	150	-	✓	✓			
23135-22-0	Oxamyl	μg/L	50	-	15	0	<20	-	19	0	<2	-		,	/		
131-57-7	Oxybenzone	μg/L	-	-	3	3	0.032	-	13	11	0.029	-	✓	✓			







						Producti	on Wells			Monitor	ing Wells			Meets	COPC Criter	Screeni ia*	ing
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Initia 2		Secoi 5	ndary 6
95-47-6	o-Xylene	μg/L	-	-	201	0	<0.5	-	475	1	0.31	-					
4685-14-7	Paraquat	μg/L	-	-	3	0	<13	-	19	0	<32	-					
189084-64-8	PBDE-100	μg/L	-	-	3	0	<0.005	-	13	0	<0.05	-					
182677-30-1	PBDE-138	μg/L	-	-	3	0	<0.005	-	13	0	<0.05	-					
68631-49-2	PBDE-153	μg/L	-	-	3	0	<0.005	-	13	0	<0.05	-					
207122-15-4	PBDE-154	μg/L	-	-	3	0	<0.005	-	13	0	<0.05	-					
147217-75-2	PBDE-17	μg/L	-	-	3	0	<0.005	-	13	0	<0.05	-					
41318-75-6	PBDE-28	μg/L	-	-	3	0	<0.005	-	13	0	< 0.05	-					
5436-43-1	PBDE-47	μg/L	-	-	3	0	<0.005	-	13	0	<0.05	-					
243982-82-3	PBDE-49	μg/L	-	-	3	0	< 0.005	-	13	0	< 0.05	-					
182346-21-0	PBDE-85	μg/L	-	-	3	0	<0.005	-	13	0	<0.05	-					
60348-60-9	PBDE-99	μg/L	-	-	3	0	<0.005	-	13	0	< 0.05	-					
76-01-7	Pentachloroethane	μg/L	-	-	4	0	<5	-	20	0	<5	-					
82-68-8	Pentachloronitrobenzene	μg/L	-	-	3	0	<1	-	19	0	<1	-					
87-86-5	Pentachlorophenol	μg/L	1	-	21	0	<10	-	57	2	0.53	0.53		~			
110-62-3	Pentanal	μg/L	-	-	3	0	<2	-	14	0	<2	-					
14797-73-0	Perchlorate	μg/L	6	-	9	1	1.6	0.27	19	8	6.7	1.12	✓	✓ v	<b>✓</b>	✓	
375-73-5	PFBS	μg/L	-	0.5	8	6	0.0051	0.01	16	11	0.0058	0.01	✓	✓ v	′		
335-76-2	PFDA	μg/L	-	-	8	2	0.0017	-	16	6	0.0062	-	✓	✓			
307-55-1	PFDoA	μg/L	-	-	8	2	0.0017	-	16	5	0.0018	-	✓	✓			
375-85-9	PFHpA	μg/L	-	-	8	8	0.0027	-	16	16	0.0055	-	✓	✓			
307-24-4	PFHxA	μg/L	-	-	8	6	0.012	-	16	14	0.013	-	✓	✓			
355-46-4	PFHxS	μg/L	-	-	8	6	0.0048	-	16	11	0.0065	-	✓	✓			
375-95-1	PFNA	μg/L	-	-	8	3	0.0017	-	16	4	0.0047	-	✓	✓			
335-67-1	PFOA	μg/L	-	0.005	8	6	0.0029	0.57	16	11	0.0033	0.65	✓	✓ v	<b>✓</b>	✓	✓
1763-23-1	PFOS	μg/L	-	0.007	8	2	0.0017	0.26	16	1	0.0042	0.65	✓	٧	<b>✓</b>	✓	
376-06-7	PFTeDA	μg/L	-	-	6	0	<0.002	-	16	0	<0.002	-					
72629-94-8	PFTrDA	μg/L	-	-	8	2	0.0017	-	16	1	0.00041	-	✓				
2058-94-8	PFUnA	μg/L	-	-	8	2	0.0017	_	16	8	0.0031	_	<b>√</b>	✓			
TPHd	PHC AS DIESEL FUEL	μg/L	-	-	0	0	-	_	3	1	100	_					
85-01-8	Phenanthrene	μg/L	-	-	6	0	<5	_	38	0	<10	_					
108-95-2	Phenol	μg/L	_	_	3	0	<1	_	19	1	1.2	_					
57-41-0	Phenytoin (Dilantin)	μg/L	_	_	3	0	<0.06	_	13	0	<0.005	_					
7723-14-0	Phosphorus	μg/L	_	_	3	3	620	_	13	13	730	_	<b>✓</b>	<b>√</b>			
85-44-9	Phthalic Anhydride	μg/L μg/L	_	_	0	0	-	_	5	5	15	_		<b>√</b>			
1918-02-1	Picloram	μg/L μg/L	500	_	11	0	<1	_	19	0	<0.6	_		_	_		
99-87-6	p-Isopropyltoluene	μg/L μg/L	-	_	34	28	0.5	_	475	0	<1	_	<b>√</b>				
7440-09-7	Potassium	μg/L μg/L	_	_	12	12	3900	_	32	32	12000	_	<b>√</b>	<b>√</b>			
55268-74-1	Praziquantel	μg/L μg/L	_	_	3	0	<0.001	_	13	1	0.0066	_					
125-33-7	Primidone	μg/L μg/L	_	_	3	2	0.0034	_	13	4	0.003	_	<b>√</b>	<b>√</b>			
57-83-0	Progesterone	μg/L μg/L	_	_	3	0	<0.001	_	13	1	0.0003						$\vdash$
1610-18-0	Prometon	μg/L μg/L		_	5	0	<1	_	19	0	<1	_					$\vdash$
7287-19-6	Prometryn	μg/L μg/L	_	_	5	0	<2	_	19	0	<1	-					$\vdash$
	Propachlor (2-Chloro-n-(1-Methylethy)-	μу/∟	-		J	0	~2	_	13	0		-					$\vdash$
1918-16-7	n-phenylac	μg/L	-	90	8	0	<2	-	38	0	<2.5	-		·			







						Production Wells Monitoring Wells						Meets COPC Screen Criteria*						
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Initio	al 3 4	Seco.	ndary 6	
123-38-6	Propanal	μg/L	-	-	3	0	<2	-	14	2	1.4	-						
114-26-1	Propoxur (Baygon)	μg/L	-	-	5	0	<10	-	19	0	<2	-						
129-00-0	Pyrene	μg/L	-	-	6	0	<5	-	38	0	<10	-						
110-86-1	Pyridine	μg/L	-	-	3	0	<5	-	19	0	<50	-						
91-22-5	Quinoline	μg/L	-	-	3	0	< 0.002	-	13	10	0.006	-		✓				
13982-63-3	Radium 226	pci/L	-	-	6	3	0.643	-	19	19	0.523	=	✓	✓				
15262-20-1	Radium 228	pci/L	-	-	6	3	0.683	-	19	19	0.998	-	✓	✓				
121-82-4	RDX	μg/L	-	0.3	3	0	<5	-	13	0	<5	-		,	✓			
69-72-7	Salicylic Acid	μg/L	-	-	3	3	1	-	13	13	0.74	-	✓	✓				
135-98-8	sec-Butylbenzene	μg/L	-	260	34	0	<0.5	-	475	0	<1	-			✓			
7782-49-2	Selenium	μg/L	50	-	12	10	4.6	0.09	38	34	9	0.18	✓	✓ .	✓ ✓			
7631-86-9	Silica	μg/L	-	-	9	9	50000	-	26	26	44000	-	<b>✓</b>	✓				
7440-22-4	Silver	μg/L	-	-	12	0	<0.2	-	38	0	<0.2	-						
122-34-9	Simazine	μg/L	4	-	15	0	<1	-	19	0	<1	-		,	✓			
7440-23-5	Sodium	μg/L	-	-	12	12	96000	-	32	32	120000	-	✓	✓				
10098-97-2	Strontium-90	pci/L	8	-	3	3	0.255	0.03	19	19	0.461	0.06	✓	✓ ,	✓ ✓			
100-42-5	Styrene	μg/L	100	-	34	0	<0.5	-	475	0	<1	-			✓			
56038-13-2	Sucralose	μg/L	-	-	3	2	0.34	-	13	1	0.0086	-	✓					
723-46-6	Sulfamethoxazole	μg/L	-	-	3	2	0.00026	-	13	1	0.00027	-	✓					
14808-79-8	Sulfate as SO4	μg/L	-	-	9	9	346000	-	24	24	1600000	-	✓	✓				
18496-25-8	Sulfide	μg/L	-	-	3	0	<100	-	13	0	<100	-						
115-96-8	TCEP	μg/L	-	-	3	1	0.00054	-	13	3	0.002	-	✓	✓				
13674-84-5	ТСРР	μg/L	-	-	3	2	0.041	-	13	11	0.008	-	✓	✓				
13674-87-8	TDCPP	μg/L	-	-	3	2	0.0034	-	13	11	0.0041	-	✓	✓				
TIC	Tentatively Identified Compounds	μg/L	-	-	9	3	9.7	-	60	35	43	-	✓	✓				
5902-51-2	Terbacil	μg/L	-	-	5	0	<20	-	19	0	<20	-						
13071-79-9	Terbufos	μg/L	-	-	3	0	<0.03	-	13	0	< 0.03	-						
98-06-6	tert-Butylbenzene	μg/L	-	260	34	0	<0.5	-	475	0	<1	-		,	<b>✓</b>			
994-05-8	Tertiary-amyl methyl ether	μg/L	-	_	197	0	<3	-	454	0	<1	-						
75-65-0	Tertiary-butyl alcohol	μg/L	-	12	103	0	<2	-	467	1	8.1	0.68		,	<b>√</b>			
58-22-0	Testosterone	μg/L	-	-	3	0	<0.001	-	13	3	0.00034	-		✓				
127-18-4	Tetrachloroethene	μg/L	5	-	207	205	54	10.80	476	288	470	94.00	✓	✓ ,	<b>√</b>	✓	✓	
109-99-9	Tetrahydrofuran	μg/L	-	-	4	0	<5	-	20	0	<5	-						
479-45-8	Tetryl	μg/L	-	-	3	0	<5	-	13	0	<5	-						
7440-28-0	Thallium	μg/L	2	-	12	1	0.033	0.02	38	5	0.05	0.03	✓	✓ ,	<b>√</b>			
28249-77-6	Thiobencarb	μg/L	70	-	15	0	<1	-	19	0	<1	-		٠,	✓			
7440-29-1	Thorium	μg/L	-	-	6	2	0.42	-	26	10	0.75	-	✓	✓				
ODOR	Threshold Odor Number	Units	-	-	9	3	100	-	19	19	800	-	✓	✓				
108-88-3	Toluene	μg/L	150	-	199	1	0.6	0.00	475	11	1.6	0.01	✓	✓ ,	<b>√</b>			
TDS	Total Dissolved Solids	μg/L	-	-	9	9	1232000	-	24	24	1970000	-	✓	✓				
TOC	Total Organic Carbon	μg/L	-	-	3	3	1400	-	13	13	21000	-	✓	✓				
1336-36-3	Total PCBs	μg/L	0.5	-	11	0	<0.5	_	19	0	<25	_		,	✓			
TSuS	Total Suspended Solids	μg/L	-	-	3	3	54000	_	13	13	150000	_	<b>√</b>	✓				
THM	Total Trihalomethanes	μg/L	80	-	34	33	4.8	0.06	19	13	18	0.23	<b>√</b>	✓ ,	✓ ✓			
1330-20-7	Total Xylenes	μg/L	1750	<del>                                     </del>	34	0	<0.5		19	0	<0.5	-			<b>√</b>			







						Producti	on Wells			Monitor	ing Wells		I		COPC Criter	Screen ia*	iing
CASRN	Parameter	Units	MCL	NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	Number of Observations	Number of Detections	Maximum Observed Value	Ratio of Maximum to MCL or NL	1	Initial 2 3	4	Seco 5	ondary 6
8001-35-2	Toxaphene	μg/L	3	-	15	0	<1	-	19	0	<50	-		✓			
TPHg	TPHg	μg/L	-	-	0	0	-	-	45	0	<50	-					
156-60-5	trans-1,2-Dichloroethene	μg/L	10	-	201	0	<0.5	-	475	8	15	1.50		✓ ✓	✓	✓	✓
10061-02-6	trans-1,3-Dichloropropene	μg/L	-	-	34	0	<0.5	-	475	0	<1	-					
110-57-6	trans-1,4-Dichloro-2-butene	μg/L	-	-	4	0	<0.5	-	20	0	<0.5	-					
24017-47-8	Triazofos	μg/L	-	-	4	0	<2	-	19	0	<2	-					
75-96-7	Tribromoacetic acid	μg/L	-	-	0	0	-	-	6	0	<2	-					
76-03-9	Trichloroacetic acid	μg/L	-	ı	3	0	<1	-	19	1	0.87	-					
79-01-6	Trichloroethene	μg/L	5	ı	205	200	63.2	12.64	475	227	370	74.00	✓	✓ ✓	✓	✓	✓
75-69-4	Trichlorofluoromethane	μg/L	150	ı	34	0	<5	-	475	4	1.6	0.01		✓ ✓			
3380-34-5	Triclosan	μg/L	-	ı	3	1	0.0093	-	13	0	< 0.002	-	✓				
1582-09-8	Trifluralin	μg/L	-	-	8	0	<1	-	38	0	<1	-					
738-70-5	Trimethoprim	μg/L	-	-	3	2	0.00054	-	13	2	0.00048	-	✓				
786-19-6	Trithion	μg/L	-	ı	3	0	<1	-	19	0	<1	-					
10028-17-8	Tritium	pci/L	20000	ı	3	3	57.5	0.00	19	19	70	0.00	✓	✓ ✓			
TURB	Turbidity	NTU	-	-	9	9	48	-	19	19	38	-	✓	✓			
7440-61-1	Uranium Rad	pci/L	20	ı	7	7	14	0.70	19	19	14	0.70	✓	✓ ✓	✓	✓	
7440-62-2	Vanadium	μg/L	-	50	10	10	14	0.28	26	25	11	0.22	✓	✓ ✓	✓		
108-05-4	Vinyl Acetate	μg/L	-	-	4	0	<5	-	20	0	<5	-					
75-01-4	Vinyl Chloride	μg/L	0.5	1	34	0	<0.5	-	475	12	32	64.00		✓ ✓	<b>√</b>	✓	✓
7440-66-6	Zinc	μg/L	-	-	12	6	110	-	38	32	23	-	✓	<b>√</b>			

#### Notes:

CASRN = Chemical Abstracts Service Registry Number (as applicable); µg/L = micrograms per liter; ng/L = nanograms per liter; mpl/ 100mL = most probable number per 100 milliliters; NTU = nephelometric turbidity units;

\*Initial and Secondary COPC screening criteria are described in Section 5 of the report, as follows:

#### *Initial screening criteria:*

- 1. Constituents that were detected in the four production wells (SM-3, SM-4, SM-8, or SM-9);
- 2. Constituents that were non-detect in the four production wells, but had three or more detections in one or more monitoring wells;
- 3. Detected constituents with a primary maximum contaminant level (MCL) or notification level (NL); and,
- 4. Constituents meeting screening criteria 1 or 2 and screening criteria 3 with a ratio of detected concentration (maximum observed) to MCL or NL greater than 0.05 (5%).

#### Secondary screening criteria:

- 5. Constituents meeting initial screening criteria with a ratio of detected concentration (maximum observed) to MCL or NL greater than 0.5 (50%); and,
- 6. Constituents meeting initial screening criteria and screening criteria 5 which are synthetic organics (these constituents are most likely to drive design for the planned treatment system using ultraviolet/advanced oxidation process and granular activated carbon).

Full Raw Water Quality Characterization - Olympic Well Field - Step 2 of 97-005 Evaluation Current Rev No.: 1: 308038-03533-00-EN-REP-0004







## Appendix C Tentatively Identified Compound (TIC) Results







	Modeled Capture							
Well ID	Zone	Sampling Date	Analytical Method	CASRN	Chemical Name	Result	Units	Comments
SM-9	SM-9	2020-05-27	E625.1	7642-04-8	2-Octene, (Z)-	34	μg/L	tic_retention_time 1.582, 2-Octene, (Z)-
SM-9	SM-9	2020-05-27	E625.1	TIC	Tentatively Identified Compounds	9.7	μg/L	Unknown #1 (possible alkane MW=140)
SM-9	SM-9	2020-05-27	E625.1	TIC	Tentatively Identified Compounds	8.8	μg/L	Unknown #2 (possible alkane MW=158)
SM-9	SM-9	2020-05-27	E525.2	TIC	Tentatively Identified Compounds	2.8	μg/L	Unknown #1 (possible ester MW=368)
SM-9	SM-9	2020-05-27	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
SM-9	SM-9	2020-05-27	E625.1	127-18-4	Tetrachloroethene	12	μg/L	Tetrachloroethene
SM-8	SM-8	2020-06-09	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
SM-8	SM-8	2020-06-09	E525.2	TIC	Tentatively Identified Compounds	0	μg/L	
SM-8	SM-8	2020-06-09	E625.1	TIC	Tentatively Identified Compounds	0	μg/L	
SM-4	SM-4	2020-06-30	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
SM-4	SM-4	2020-06-30	E525.2	TIC	Tentatively Identified Compounds	0	μg/L	
SM-4	SM-4	2020-06-30	E625.1	127-18-4	Tetrachloroethene	13	μg/L	Tetrachloroethene
OB-7	SM-4	2020-06-03	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-7	SM-4	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-7	SM-4	2020-06-03	E625.1	TIC	Tentatively Identified Compounds	0	μg/L	
OB-5	SM-4	2020-06-05	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-5	SM-4	2020-06-05	E525.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-5	SM-4	2020-06-05	E625.1	TIC	Tentatively Identified Compounds	0	μg/L	
OB-3	SM-8	2020-06-03	E525.2	934-34-9	2(3H)-Benzothiazolone	1.4	μg/L	tic_retention_time 7.991, 2(3H)-Benzothiazolone
OB-3	SM-8	2020-06-03	E625.1	95-16-9	Benzothiazole	8.5	μg/L	tic_retention_time 7.452, Benzothiazole
OB-3	SM-8	2020-06-03	E525.2	95-16-9	Benzothiazole	4	μg/L	tic_retention_time 5.122, Benzothiazole
OB-3	SM-8	2020-06-03	E525.2	1122-82-3	Cyclohexane, isothiocyanato-	1	μg/L	tic_retention_time 5.172, Cyclohexane, isothiocyanato-
OB-3	SM-8	2020-06-03	E525.2	85-44-9	Phthalic Anhydride	15	μg/L	tic_retention_time 5.726, Phthalic Anhydride
OB-3	SM-8	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	26	μg/L	Unknown #8 (possible ketone MW=250); tic_retention_time 10.336
OB-3	SM-8	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	20	μg/L	Unknown #4 (possible alcohol MW=178); tic_retention_time 9.591
OB-3	SM-8	2020-06-03	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-3	SM-8	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	9.8	μg/L	Unknown #6 (possible amide MW=205); tic_retention_time 9.863
OB-3	SM-8	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	6.7	μg/L	Unknown #9 (possible ketone MW=262); tic_retention_time 11.293
OB-3	SM-8	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	3.3	μg/L	Unknown #5 (possible aldehyde MW=234); tic_retention_time 9.722
OB-3	SM-8	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	1.9	μg/L	Unknown #1 (possible nitrile MW=191); tic_retention_time 8.816
OB-3	SM-8	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	1.5	μg/L	Unknown #2 (possible hydrazone MW=100); tic_retention_time 9.018
OB-3	SM-8	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	1.2	μg/L	Unknown #3 (possible PAH MW=234); tic_retention_time 9.279
OB-3	SM-8	2020-06-03	E525.2	TIC	Tentatively Identified Compounds	1	μg/L	Unknown #7 (possible alcohol MW=164); tic_retention_time 10.075
OB-2	SM-8	2020-06-04	E525.2	85-44-9	Phthalic Anhydride	2.2	μg/L	tic_retention_time 5.716, Phthalic Anhydride
OB-2	SM-8	2020-06-04	E625.1	TIC	Tentatively Identified Compounds	43	μg/L	Unknown #1 (possible alcohol MW=130); tic_retention_time 5.587
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	18	μg/L	Unknown #2 (possible alcohol MW=144); tic_retention_time 6.129
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	15	μg/L	Unknown #15 (possible alkene MW=262); tic_retention_time 10.658
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	13	μg/L	Unknown #12 (possible alkane MW=248); tic_retention_time 9.863







	Modeled Capture							
Well ID	Zone	Sampling Date	Analytical Method	CASRN	Chemical Name	Result	Units	Comments
OB-2	SM-8	2020-06-04	E625.1	TIC	Tentatively Identified Compounds	13	μg/L	Unknown #3 (possible diazine MW=205); tic_retention_time 12.42
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	11	μg/L	Unknown #6 (possible PAH MW=188); tic_retention_time 8.474
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	8.4	μg/L	Unknown #11 (possible phenol MW=178); tic_retention_time 9.581
OB-2	SM-8	2020-06-04	E625.1	TIC	Tentatively Identified Compounds	8.1	μg/L	Unknown #2 (possible alkane MW=126); tic_retention_time 11.444
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	6.3	μg/L	Unknown #13 (possible amine MW=212); tic_retention_time 10.175
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	5.8	μg/L	Unknown #8 (possible amine MW=129); tic_retention_time 9.007
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	5.2	μg/L	Unknown #7 (possible ether MW=148); tic_retention_time 8.967
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	4.8	μg/L	Unknown #1 (possible carboxylic acid MW=242); tic_retention_time 5.676
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	4.1	μg/L	Unknown #4 (possible carboxylic acid MW=284); tic_retention_time 7.719
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	4	μg/L	Unknown #16 (possible ketone MW=260); tic_retention_time 11.222
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	2.9	μg/L	Unknown #14 (possible pyridine MW=180); tic_retention_time 10.568
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	2	μg/L	Unknown #9 (possible carboxylic acid MW=234); tic_retention_time 9.209
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	1.8	μg/L	Unknown #10 (possible phenol MW=164); tic_retention_time 9.279
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	1.8	μg/L	Unknown #5 (possible amide MW=175); tic_retention_time 8.293
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	1.8	μg/L	Unknown #17 (possible azole MW=178); tic_retention_time 11.363
OB-2	SM-8	2020-06-04	E525.2	TIC	Tentatively Identified Compounds	1.7	μg/L	Unknown #3 (possible carboxylic acid MW=268); tic_retention_time 6.35
OB-2	SM-8	2020-06-04	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-17C	SM-4	2020-05-27	E525.2	TIC	Tentatively Identified Compounds	4.6	μg/L	Unknown #2 (possible ketone MW=140); tic_retention_time 10.668
OB-17C	SM-4	2020-05-27	E525.2	TIC	Tentatively Identified Compounds	1.5	μg/L	Unknown #1 (possible alkyne MW=160); tic_retention_time 8.816
OB-17C	SM-4	2020-05-27	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-17C	SM-4	2020-05-27	E625.1	TIC	Tentatively Identified Compounds	0	μg/L	
OB-17B	SM-4	2020-05-27	E625.1	7642-04-8	2-Octene, (Z)-	30	μg/L	tic_retention_time 1.588, 2-Octene, (Z)-
OB-17B	SM-4	2020-05-27	E525.2	85-44-9	Phthalic Anhydride	1.3	μg/L	tic_retention_time 5.766, Phthalic Anhydride
OB-17B	SM-4	2020-05-27	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-14C	SM-9	2020-06-02	E625.1	7642-04-8	2-Octene, (Z)-	27	μg/L	tic_retention_time 1.582. 2-Octene, (Z)-
OB-14C	SM-9	2020-06-02	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-14C	SM-9	2020-06-02	E525.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-14B	SM-9	2020-06-02	E625.1	7642-04-8	2-Octene, (Z)-	250	μg/L	tic_retention_time 1.582. 2-Octene, (Z)-
OB-14B	SM-9	2020-06-02	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-14B	SM-9	2020-06-02	E525.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-13C	SM-9	2020-05-29	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-13C	SM-9	2020-05-29	E525.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-13C	SM-9	2020-05-29	E625.1	TIC	Tentatively Identified Compounds	0	μg/L	
OB-12C	SM-4	2020-05-26	E525.2	85-44-9	Phthalic Anhydride	4.6	μg/L	00:05:46, Phthalic Anhydride
OB-12C	SM-4	2020-05-26	E525.2	TIC	Tentatively Identified Compounds	6.7	μg/L	00:10:41
OB-12C	SM-4	2020-05-26	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-12C	SM-4	2020-05-26	E625.1	TIC	Tentatively Identified Compounds	0	μg/L	
OB-12B	SM-4	2020-05-26	E625.1	7642-04-8	2-Octene, (Z)-	26	μg/L	00:01:35, 2-Octene, (Z)-







Well ID	Modeled Capture Zone	Sampling Date	Analytical Method	CASRN	Chemical Name	Result	Units	Comments
OB-12B	SM-4	2020-05-26	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-12B	SM-4	2020-05-26	E525.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-11C	SM-4	2020-05-28	E625.1	111-66-0	Octene-1	25	μg/L	tic_retention time 1.575, Octene-1
OB-11C	SM-4	2020-05-28	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-11C	SM-4	2020-05-28	E525.2	TIC	Tentatively Identified Compounds	0	μg/L	
OB-11C	SM-4	2020-05-28	E625.1	127-18-4	Tetrachloroethene	55	μg/L	tic_retention time 1.699, Tetrachloroethene
OB-11B	SM-4	2020-05-28	E625.1	7642-04-8	2-Octene, (Z)-	30	μg/L	tic_retention time 1.576, 2-Octene, (Z)-
OB-11B	SM-4	2020-05-28	E525.2	85-44-9	Phthalic Anhydride	8.3	μg/L	tic_retention time 5.776, Phthalic Anhydride
OB-11B	SM-4	2020-05-28	E525.2	TIC	Tentatively Identified Compounds	2.4	μg/L	possible ester MW=274; tic_retention time 7.84
OB-11B	SM-4	2020-05-28	E525.2	TIC	Tentatively Identified Compounds	1.4	μg/L	possible nitrile MW=191; tic_retention time 8.857
OB-11B	SM-4	2020-05-28	E525.2	TIC	Tentatively Identified Compounds	1.4	μg/L	possible diazine MW=205; tic_retention time 9.893
OB-11B	SM-4	2020-05-28	E524.2	TIC	Tentatively Identified Compounds	0	μg/L	

#### Notes:

CASRN = Chemical Abstracts Service Registry Number (as applicable);  $\mu g/L = micrograms$  per liter; '--' = not applicable.







### Appendix D Microbiological Quality Data







Well ID	Modeled Capture Zone	Sampling Date	Method	Chemical Name	Result	Units	RL
SM-4	SM-4	2020-06-30	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
SM-8	SM-8	2020-06-09	A9221E	E.Coli	< 1.1	mpn/100ml	1.1
SM-9	SM-9	2020-05-27	A9215B	E.Coli	< 1.1	mpn/100ml	1.1
OB-11B	SM-4	2020-05-28	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-11C	SM-4	2020-05-28	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-12B	SM-4	2020-05-26	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-12C	SM-4	2020-05-26	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-13C	SM-9	2020-05-29	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-14B	SM-9	2020-06-02	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-14C	SM-9	2020-06-02	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-17B	SM-4	2020-05-27	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-17C	SM-4	2020-05-27	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-2	SM-8	2020-06-04	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-3	SM-8	2020-06-03	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-5	SM-4	2020-06-05	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
OB-7	SM-4	2020-06-03	A9221F	E.Coli	< 1.1	mpn/100ml	1.1
SM-4	SM-4	2020-06-30	A9221B	Fecal coliform	< 1.1	mpn/100ml	1.1
SM-8	SM-8	2020-06-09	A9221F	Fecal coliform	< 1.1	mpn/100ml	1.1
SM-9	SM-9	2020-05-27	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-11B	SM-4	2020-05-28	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-11C	SM-4	2020-05-28	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-12B	SM-4	2020-05-26	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-12C	SM-4	2020-05-26	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-13C	SM-9	2020-05-29	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-14B	SM-9	2020-06-02	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-14C	SM-9	2020-06-02	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-15C	SM-9	2020-06-09	A9221E	Fecal coliform	2.2	mpn/100ml	1.1
OB-17B	SM-4	2020-05-27	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-17C	SM-4	2020-05-27	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-2	SM-8	2020-06-04	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-3	SM-8	2020-06-03	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-5	SM-4	2020-06-05	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-6D	SM-4	2020-06-10	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-6D	SM-4	2020-06-10	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
OB-7	SM-4	2020-06-03	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
SMB1-#4	SM-9	2020-06-11	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
SMB1-#4	SM-9	2020-09-03	A9222D	Fecal coliform	< 2.0	cfu/100ml	2.0
SMB1-B#1	SM-9	2020-06-11	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
SMB1-B#1	SM-9	2020-06-11	A9221E	Fecal coliform	< 1.1	mpn/100ml	1.1
SMB1-B#1	SM-9	2020-09-03	A9222D	Fecal coliform	8.0	cfu/100ml	2.0
SM-4	SM-4	2020-06-30	A9215B	Heterotrophic Plate Count	22	cfu/ml	1.0
SM-8	SM-8	2020-06-09	A9221B	Heterotrophic Plate Count	360	cfu/ml	1.0
SM-9	SM-9	2020-05-27	A9221F	Heterotrophic Plate Count	400	cfu/ml	1.0







Well ID	Modeled Capture Zone	Sampling Date	Method	Chemical Name	Result	Units	RL
OB-11B	SM-4	2020-05-28	A9215B	Heterotrophic Plate Count	230	cfu/ml	1.0
OB-11C	SM-4	2020-05-28	A9215B	Heterotrophic Plate Count	400	cfu/ml	1.0
OB-12B	SM-4	2020-05-26	A9215B	Heterotrophic Plate Count	40	cfu/ml	1.0
OB-12C	SM-4	2020-05-26	A9215B	Heterotrophic Plate Count	< 1.0	cfu/ml	1.0
OB-13C	SM-9	2020-05-29	A9215B	Heterotrophic Plate Count	140	cfu/ml	1.0
OB-14B	SM-9	2020-06-02	A9215B	Heterotrophic Plate Count	220	cfu/ml	1.0
OB-14C	SM-9	2020-06-02	A9215B	Heterotrophic Plate Count	160	cfu/ml	1.0
OB-17B	SM-4	2020-05-27	A9215B	Heterotrophic Plate Count	2.0	cfu/ml	1.0
OB-17C	SM-4	2020-05-27	A9215B	Heterotrophic Plate Count	120	cfu/ml	1.0
OB-2	SM-8	2020-06-04	A9215B	Heterotrophic Plate Count	44	cfu/ml	1.0
OB-3	SM-8	2020-06-03	A9215B	Heterotrophic Plate Count	40	cfu/ml	1.0
OB-5	SM-4	2020-06-05	A9215B	Heterotrophic Plate Count	41	cfu/ml	1.0
OB-7	SM-4	2020-06-03	A9215B	Heterotrophic Plate Count	6.0	cfu/ml	1.0
SM-4	SM-4	2020-06-30	A9221E	Total coliform	6.9	mpn/100ml	1.1
SM-8	SM-8	2020-06-09	A9215B	Total coliform	< 1.1	mpn/100ml	1.1
SM-9	SM-9	2020-05-27	A9221B	Total coliform	2.2	mpn/100ml	1.1
OB-11B	SM-4	2020-05-28	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-11C	SM-4	2020-05-28	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-12B	SM-4	2020-05-26	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-12C	SM-4	2020-05-26	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-13C	SM-9	2020-05-29	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-14B	SM-9	2020-06-02	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-14C	SM-9	2020-06-02	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-15C	SM-9	2020-06-09	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-17B	SM-4	2020-05-27	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-17C	SM-4	2020-05-27	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-2	SM-8	2020-06-04	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-3	SM-8	2020-06-03	A9221B	Total coliform	1.1	mpn/100ml	1.1
OB-5	SM-4	2020-06-05	A9221B	Total coliform	1.1	mpn/100ml	1.1
OB-6D	SM-4	2020-06-10	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-6D	SM-4	2020-06-10	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
OB-7	SM-4	2020-06-03	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
SMB1-#4	SM-9	2020-06-11	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
SMB1-#4	SM-9	2020-09-03	A9222B	Total coliform	30	cfu/100ml	2.0
SMB1-B#1	SM-9	2020-06-11	A9221B	Total coliform	1.1	mpn/100ml	1.1
SMB1-B#1	SM-9	2020-06-11	A9221B	Total coliform	< 1.1	mpn/100ml	1.1
SMB1-B#1	SM-9	2020-09-03	A9222B	Total coliform	110	cfu/100ml	2.0

#### Notes:

cfu = colony forming units; mpn = most probable number; ml = milliliter; RL = Reporting limit







# Appendix E Statistical Analysis Results and Production Well Concentrations Estimate







				Us	sing UCL95 f	or Monitoring \	Well Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo	or Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
2567.62.2	4 (2 4 5)	C-Zone	ND	ND	ND	1	ND	NB	ND	ND	ND	1	ND	NB
3567-62-2	1-(3,4-Dichlorophenyl)-3-methylurea [µg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
2327-02-8	1-(3,4-Dichlorophenyl)urea [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
2327-02-0	1-(3,4-Dichiorophenyi)urea [µg/L]	D-Zone	NA	NA	NA	שוו	ND	ND	NA	NA	NA	שוא	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
630-20-6	1,1,1,2-Tetrachloroethane [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
030-20-0	1,1,1,2-Tetrachioroethane [µg/L]	D-Zone	ND	ND	ND	IND	IND	IND	ND	ND	ND	I ND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
71-55-6	1,1,1-Trichloroethane [µg/L]	C-Zone	ND	ND	ND	ND ND	ND	ND	ND	ND	ND	ND	ND	ND
/1-33-0	1,1,1-Themoroethane [μg/L]	D-Zone	ND	ND	ND	שוו	IND	IND	ND	ND	ND	] ND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
79-34-5	34-5 1,1,2,2-Tetrachloroethane [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
79-54-5	1,1,2,2-Tetrachioroethane [µg/L]	D-Zone	ND	ND	ND	שוו	IND	IND	ND	ND	ND	] ND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
79-00-5	1,1,2-Trichloroethane [µg/L]	C-Zone	1.6	ND	ND	0.27	ND	ND	1.1	ND	ND	0.18	ND	ND
/9-00-5	1,1,2-Themoroethane [µg/L]	D-Zone	ND	ND	ND	0.27	IND	IND	ND	ND	ND	0.16	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.671	-	0.15				0.517	-	0.15			
75-34-3	1,1-Dichloroethane (1,1-DCA) [µg/L]	C-Zone	0.395	0.897	0.12	0.27	0.039	0.01	0.243	0.603	0.12	0.2	0.026	0.01
75-54-5	1,1-Dichioroethane (1,1-DCA) [µg/L]	D-Zone	ND	ND	ND	0.27	0.039	0.01	ND	ND	ND	0.2	0.020	0.01
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	2.186	_	2.111				1.545	-	1.691			
75-35-4	1,1-Dichloroethene (1,1-DCE) [μg/L]	C-Zone	1.991	3.943	0.22	1.1	0.2	0.079	1.517	3.303	0.22	0.81	0.17	0.071
75-55-4	1,1-Dictiloroethene (1,1-DCE) [µg/L]	D-Zone	0.5	0.5	0.5	1.1	0.2	0.079	0.5	0.5	0.5	0.61	0.17	0.071
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
563-58-6	1,1-Dichloropropene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
303-30-0	1,1-bichioroproperie [µg/L]	D-Zone	ND	ND	ND	IND	IND	IND	ND	ND	ND	IND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
57-14-7	1,1-Dimethylhydrazine [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
37-14-7	i, i-Dimethymydiazme [μg/L]	D-Zone	NA	NA	NA	IND	IND	IND	NA	NA	NA	IND	IND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND	_			ND	-	ND			
87-61-6	1,2,3-Trichlorobenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
07-01-0	i,z,ɔ-mcmorobenzene [µg/t]	D-Zone	ND	ND	ND	ואט	ואט	IND	ND	ND	ND	שמו	שוו	שוו
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Us	ing UCL <u>95 f</u> o	or Monitoring V	Vell Data*			Usi	ing Mean f <u>o</u>	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concent	ation within Zone			oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
96-18-4	1,2,3-Trichloropropane (1,2,3-TCP) [µg/L]	C-Zone	0.0051	0.0156	ND	0.037	0.015	0.014	0.0048	0.012	ND	0.037	0.015	0.014
90-10-4	1,2,3-ПСПЮГОРГОРАПЕ (1,2,3-ТСР) [µg/L]	D-Zone	0.19	0.19	0.19	0.057	0.015	0.014	0.19	0.19	0.19	0.057	0.015	0.014
		Sunnyside	0.0041	0.0041	0.0041				0.0041	0.0041	0.0041			
		B-Zone	ND	-	ND				ND	-	ND			
120-82-1	1,2,4-Trichlorobenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
.20 02 .	., <u>-,</u>	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
95-63-6	1,2,4-Trimethylbenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	,	D-Zone	ND	ND	ND				ND	ND	ND	<u> </u>		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	- ND	ND	_		
96-12-8	1,2-Dibromo-3-chloropropane [µg/L]	C-Zone D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
			ND ND	ND ND	ND				ND ND	ND ND	ND ND	_		
		Sunnyside B-Zone	ND ND	ND	ND				ND	- -	ND			
		C-Zone	ND ND	ND	ND				ND	ND	ND	1		
106-93-4	1,2-Dibromoethane [µg/L]	D-Zone	ND ND	ND	ND	ND	ND	ND	ND	ND ND	ND	ND	ND	ND
		Sunnyside	ND ND	ND	ND				ND	ND	ND			
		B-Zone	ND ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND			
95-50-1	1,2-Dichlorobenzene [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	_	ND				ND	-	ND			
		C-Zone	0.33	0.59	ND				0.33	0.47	ND	1		
107-06-2	1,2-Dichloroethane [μg/L]	D-Zone	0.15	0.15	0.15	0.083	0.034	0.0085	0.15	0.15	0.15	0.083	0.029	0.0085
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.16	-	ND				0.16	-	ND			
70.07.5	1.2 Diables	C-Zone	ND	ND	ND	0.040	ND	ND	ND	ND	ND	0.048	ND	ND
78-87-5	1,2-Dichloropropane [µg/L]	D-Zone	ND	ND	ND	0.048	ND	ND	ND	ND	ND	0.048	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
122-66-7	1,2-Diphenylhydrazine [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
122-00-7	r,z-σιρπεπушуαrazme [μg/ t]	D-Zone	ND	ND	ND	ND	IND	IND	ND	ND	ND	IND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
108-67-8	1,3,5-Trimethylbenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
100 07 0	1,5,5 11εαιγισείτ2ειτε [μg/ 2]	D-Zone	ND	ND	ND	115			ND	ND	ND		1,15	1,15
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
99-35-4	1,3,5-Trinitrobenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	, , ,	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
544.70.4	42 P: 11 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	C-Zone	ND	ND	ND		NB	ND	ND	ND	ND	1	ND	NB
541-73-1	1,3-Dichlorobenzene [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
142-28-9	12 Dichlerencenene [ug/l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
142-20-9	1,3-Dichloropropane [µg/L]	D-Zone	ND	ND	ND	IND	ND	ND	ND	ND	ND	] ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
542-75-6	1,3-Dichloropropene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
342-73-0	1,3-ыстогоргорене [µg/с]	D-Zone	ND	ND	ND	ND	IND	ND	ND	ND	ND	IND	IND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
99-65-0	1,3-Dinitrobenzene [µg/L]	C-Zone	0.26	ND	ND	0.044	ND	ND	0.26	ND	ND	0.044	ND	ND
33 03 0	1,5 Dillittoberizerie [µg/L]	D-Zone	NA	NA	NA	0.044	ND	ND	NA	NA	NA	0.044	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
106-46-7	1,4-Dichlorobenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
100 40 7	i, a biemorobenzene [pg/L]	D-Zone	ND	ND	ND		110	110	ND	ND	ND		110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	102.1	-	117.4	-			56.3	-	106.6			
123-91-1	1,4-Dioxane [µg/L]	C-Zone	27.83	48.38	1.787	35.4	2.1	2.2	22.7	36.32	2.298	20.8	1.6	2
.25 51 1	.,	D-Zone	ND	ND	ND	]			ND	ND	ND			_
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	-			ND	-	ND	_		
763051-92-9	11Cl-PF3OUdS [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND	-			ND	ND	ND	_		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND	-			ND	-	ND	1		
57-91-0	17-a-Estradiol [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	5	D-Zone	NA	NA	NA	-			NA	NA	NA	-		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	- ND	ND	-			ND	- ND	ND	-		
57-63-6	17-a-Ethynylestradiol [µg/L]	C-Zone	ND NA	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	NA NA	NA	NA	-			NA NA	NA	NA	-		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	- ND	ND	-			ND ND	- ND	ND	-		
50-28-2	17-b-Estradiol [μg/L]	C-Zone	ND NA	ND NA	ND NA	ND	ND	ND	ND NA	ND NA	ND NA	ND	ND	ND
		D-Zone	NA NA	NA NA	NA NA	-			NA NA	NA NA	NA NA	-		
		Sunnyside	NA	NA					NA ND	NA	NA			
		B-Zone	ND	- ND	ND	1			ND ND	- ND	ND	+		
124-19-6	1-Nonanal [µg/L]	C-Zone	ND NA	ND NA	ND NA	ND	ND	ND	ND NA		ND NA	ND	ND	ND
		D-Zone	NA NA	NA NA	NA NA	-			NA NA	NA NA	NA	-		
		Sunnyside	NA	NA	NA				NA	NA	NA			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ng Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	NA	-	NA				NA	-	NA			
024.24.0	2/211) Bannathianalana [/1]	C-Zone	NA	1.4	NA	NIA	0.061	NI A	NA	1.4	NA	NA NA	0.061	NI A
934-34-9	2(3H)-Benzothiazolone [µg/L]	D-Zone	NA	NA	NA	NA	0.061	NA	NA	NA	NA	INA	0.061	NA
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	_	ND				ND	-	ND			
594-20-7	2,2-Dichloropropane [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
394-20-1		D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	IND	IND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	_	ND				ND	-	ND			
51207-31-9	2,3,7,8-Tetrachlorodibenzofuran [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
31207-31-9	2,3,7,0-Tetrachiorodibenzordrafi [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	IND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
93-76-5	2,4,5-T [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
33 70 3	2,4,5 Γ [μg/ ε]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND		ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
93-72-1	2,4,5-TP (Silvex) [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
33 72 1	2, 1, 5 11 (5πνελ) [μβ, 2]	D-Zone	ND	ND	ND	110	110	110	ND	ND	ND		IND	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	_	ND				ND	-	ND	<u> </u>		
95-95-4	2,4,5-Trichlorophenol [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	_, ,,oee.epee. [µg, _]	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
88-06-2	2,4,6-Trichlorophenol [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	, , , , , , , , , , , , , , , , , , , ,	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND	_		
118-96-7	2,4,6-Trinitrotoluene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	NA	NA	NA				NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
94-75-7	2,4-D [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	. •	D-Zone	ND	ND	ND				ND	ND	ND	<u> </u>		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	-	ND	<u> </u>		
94-82-6	2,4-DB [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND				ND	ND	ND	<u> </u>		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	- ND	ND			
120-83-2	2,4-Dichlorophenol [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
414531			Concent	ration within	Capture	Estimated C	oncentrations fo	or Production	Concent	ration within			oncentrations f	or Production
CASRN	Constituent [Units]	Layer		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
105-67-9	2,4-Dimethylphenol [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
103-07-9	2,4-Dimethylphenol [µg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	IND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
51-28-5	2,4-Dinitrophenol [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
31-20-3	2,4-Dilliti Ophenoi [µg/L]	D-Zone	ND	ND	ND	I ND	ND	IND	ND	ND	ND	IND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
121-14-2	2,4-Dinitrotoluene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
121-14-2	2,4-Dilitiotoldene [μg/L]	D-Zone	ND	ND	ND	I ND	ND	IND	ND	ND	ND	IND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
606-20-2	2,6-Dinitrotoluene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
000-20-2	2,0-Dilitiotoldene [μg/L]	D-Zone	ND	ND	ND	I ND	ND	IND	ND	ND	ND	IND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
35572-78-2	2-Amino-4,6-Dinitrotoluene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
33372-70-2	2-Ammo-4,0-Dimitrotoidene [µg/L]	D-Zone	NA	NA	NA	I ND	ND	IND	NA	NA	NA	IND	ND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
78-93-3	2-Butanone [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
70-93-3	2-Butanone [µg/L]	D-Zone	ND	ND	ND	I ND	ND	IND	ND	ND	ND	IND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
110-75-8	2-Chloroethyl Vinyl Ether [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
110 73 0	Z Chloroethyr Villyr Ether [μg/ L]	D-Zone	ND	ND	ND	I ND	ND	ND	ND	ND	ND	I ND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
91-58-7	2-Chloronaphthalene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
31 30 7	2 emoronaphinalene [pg/2]	D-Zone	ND	ND	ND		145	110	ND	ND	ND		110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
95-57-8	2-Chlorophenol [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
33 37 0	2 (11010)11(1101 [µg/2]	D-Zone	ND	ND	ND		145	110	ND	ND	ND		110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
95-49-8	2-Chlorotoluene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	Σ σσ.σ.σ.σ.σ. [μg/ ε]	D-Zone	ND	ND	ND		. 45		ND	ND	ND	1	. 10	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND	]		
591-78-6	2-Hexanone [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
331 70 0	2 Hexanone [µg/L]	D-Zone	ND	ND	ND	140	140	IND	ND	ND	ND		ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Usi	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	r Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
91-57-6	2 Mothydnaphthalono [ug/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
91-57-0	2-Methylnaphthalene [µg/L]	D-Zone	ND	ND	ND	ND	IND	ND	ND	ND	ND		IND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
95-48-7	2-Methylphenol [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	Σ πετημετίοι [μg/ 2]	D-Zone	ND	ND	ND	113	113	115	ND	ND	ND		1,15	1,15
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
88-74-4	2-Nitroaniline [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	3· · -	D-Zone	ND	ND	ND				ND	ND	ND	<u> </u>		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	- ND	ND	_		
88-75-5	2-Nitrophenol [µg/L]	C-Zone D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
			ND ND	ND ND	ND ND				ND ND	ND ND	ND ND	_		
		Sunnyside B-Zone	ND ND	-	ND ND				ND	-	ND			
		C-Zone	ND ND	ND	ND ND				ND	ND	ND	1		
88-72-2	2-Nitrotoluene [µg/L]	D-Zone	NA NA	NA NA	NA NA	ND	ND	ND	NA NA	NA NA	NA NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA NA	NA			
		B-Zone	30	-	250				29	-	250			
		C-Zone	NA	NA	27				NA	NA	27			
7642-04-8	2-Octene, (Ζ)- [μg/L]	D-Zone	NA	NA	NA	9	NA	6	NA	NA	NA	8.6	NA	6
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND			
84989-04-8	3 & 4-Methylphenol [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	_	ND			
01.04.1	2.21 Diable as benefities [//]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
91-94-1	3,3'-Dichlorobenzidine [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	םא [	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
95-76-1	3,4-Dichloroaniline [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
93-70-1	5,4-Dictiloroarilline [µg/L]	D-Zone	NA	NA	NA	IND	IND	ND	NA	NA	NA	IND	IND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
51-36-5	3,5-Dichlorobenzoic acid [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	5,5 2.5.110.000.112010 dela [µg/ L]	D-Zone	ND	ND	ND	'''			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
16655-82-6	3-Hydroxycarbofuran [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	, , , , , , , , , , , , , , , , , , ,	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Us	ing UCL95 f	or Monitoring V	Vell Data*			Usi	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production	Concenti	ation within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
99-09-2	2 Nitropolito - Francis	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
99-09-2	3-Nitroaniline [µg/L]	D-Zone	ND	ND	ND	שא	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
99-08-1	3-Nitrotoluene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
33-00-1	3-Mitrotoldene [μg/L]	D-Zone	NA	NA	NA	I ND	ND	ND	NA	NA	NA	IND	ND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
72-54-8	4,4´-DDD [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
72 34 0	η, η υυυ [μg/ c]	D-Zone	ND	ND	ND		ND	145	ND	ND	ND	110	ND	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	_			ND	-	ND			
72-55-9	4,4´-DDE [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	,	D-Zone	ND	ND	ND	_			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	_			ND	-	ND			
50-29-3	4,4´-DDT [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND	_			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND ND	- ND	ND	-			ND ND	- ND	ND			
534-52-1	4,6-Dinitro-2-Methyl phenol [µg/L]	C-Zone D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
		Sunnyside	ND ND	ND ND	ND ND	+			ND ND	ND ND	ND			
		B-Zone	ND ND	-	ND ND				ND ND	- ND	ND			
		C-Zone	ND ND	ND	ND ND	-			ND ND	ND	ND			
19406-51-0	4-Amino-2,6-Dinitrotoluene [µg/L]	D-Zone	NA NA	NA NA	NA NA	ND	ND	ND	NA	NA NA	NA	ND	ND	ND
		Sunnyside	NA NA	NA NA	NA	1			NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND			
101-55-3	4-Bromophenyl Phenyl Ether [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND	1			ND	ND	ND			
		B-Zone	ND	-	ND				ND	_	ND			
50.50.7	4.611 2.44 1.1 1.5 (1.3	C-Zone	ND	ND	ND	1	NE		ND	ND	ND		NB	
59-50-7	4-Chloro-3-Methylphenol [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
106-47-8	4-Chloroaniline [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
100-47-0	4-Chloroaniline [µg/L]	D-Zone	ND	ND	ND	שמו	שמו	ואט	ND	ND	ND	שויו	ND	שויו
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
7005-72-3	4-Chlorophenyl Phenyl Ether [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
7003-72-3	- Chiorophenyi Frienyi Ether [µg/L]	D-Zone	ND	ND	ND		IND	ND	ND	ND	ND	IND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Us	ing UCL95 fo	or Monitoring \	Vell Data*			Usi	ng Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone			oncentrations fo Wells	r Production	Concenti	ration within Zone			oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
100.00.		C-Zone	ND	ND	ND				ND	ND	ND	1		
106-43-4	4-Chlorotoluene [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
100 10 1	A Math. I 2 and a sector of a 41	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	NB	ND	ND
108-10-1	4-Methyl-2-pentanone [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
100.01.6	A NI'. 11' F (12	C-Zone	ND	ND	ND	NE	NID	ND	ND	ND	ND		ND	ND
100-01-6	4-Nitroaniline [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
100.00.7	A NEC contract F or All	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	, NB	ND	ND
100-02-7	4-Nitrophenol [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
00.00.0	A Nitratal and For All	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	, NB	ND	ND
99-99-0	4-Nitrotoluene [μg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
104 40 5	4 Nice Johanni F. v (1)	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	, NB	ND	ND
104-40-5	4-Nonylphenol [μg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
1000 20 4	A Cost delegan al Francia	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1806-26-4	4-Octylphenol [μg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
140-66-9	A tout Ostulahanal Fug (1)	C-Zone	ND	0.014	ND	ND	0.00061	ND	ND	0.014	ND	ND	0.00061	ND
140-00-9	4-tert-Octylphenol [μg/L]	D-Zone	NA	NA	NA	שוא	0.00061	ND	NA	NA	NA	שוא	0.00061	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
2315-61-9	4-tert-Octylphenol diethoxylate [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
2515-01-9	4-tert-Octyrphenor diethoxylate [µg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA		IND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.0088	-	ND				0.0087	-	ND			
1173019-48-1	4-tert-Octylphenol monoethoxylate [µg/L]	C-Zone	ND	0.072	ND	0.0027	0.0031	ND	ND	0.052	ND	0.0026	0.0023	ND
1175019-40-1	4-tert-Octylphenol monoethoxylate [µg/L]	D-Zone	NA	NA	NA	0.0027	0.0051	ND	NA	NA	NA	0.0026	0.0023	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
756426-58-1	OCI DESCAIS [//]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
/30420-30-1	9CI-PF3ONS [μg/L]	D-Zone	ND	ND	ND	שמו	טאו	טאו	ND	ND	ND	שויו	שויו	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			







				Us	sing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
02.22.0	A 1.1 5 413	C-Zone	ND	ND	ND		NB	ND	ND	ND	ND	1	ND	N.D.
83-32-9	Acenaphthene [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
208-96-8	Acanaphthylana [ug/l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
200-90-0	Acenaphthylene [μg/L]	D-Zone	ND	ND	ND	ND	IND	ND	ND	ND	ND	שוא	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.48	-	ND				0.48	-	ND			
75-07-0	Acetaldehyde [µg/L]	C-Zone	0.37	ND	ND	0.21	ND	ND	0.37	ND	ND	0.21	ND	ND
75-07-0	Acetaideriyde [µg/L]	D-Zone	NA	NA	NA	0.21	IND	ND	NA	NA	NA	0.21	IND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
103-90-2	Acetaminophen [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
103-90-2	Acetaninophen [µg/L]	D-Zone	NA	NA	NA	I ND	IND	ND	NA	NA	NA	IND	IND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
34256-82-1	Acetochlor [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
34230 02 1	Acetochioi [µg/L]	D-Zone	ND	ND	ND	I ND	ND	IND	ND	ND	ND		ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
187022-11-3	Acetochlor ESA [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
107022 11 5	Acctocinor ESA [µg/ L]	D-Zone	NA	NA	NA	110	IND	IND	NA	NA	NA		ND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
194992-44-4	Acetochlor OA [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
134332 44 4	πεεισειποι σπ [μg/ ε]	D-Zone	NA	NA	NA		110	110	NA	NA	NA		110	110
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	20	-	ND				7.6	-	ND			
67-64-1	Acetone [µg/L]	C-Zone	31	6	ND	15	7.3	6.8	24	4.2	ND	10	7.2	6.8
		D-Zone	6.1	6.1	6.1	_			6	6	6	1		
		Sunnyside	7.4	7.4	7.4				7.4	7.4	7.4			
		B-Zone	0.62	-	ND				0.62	-	ND	1		
75-05-8	Acetonitrile [µg/L]	C-Zone	0.14	0.13	0.05	0.29	0.22	0.21	0.11	0.093	0.05	0.29	0.22	0.21
	1, 3, 3	D-Zone	ND	ND	ND				ND	ND	ND	1		
		Sunnyside	0.24	0.24	0.24				0.24	0.24	0.24			
		B-Zone	ND	-	ND				ND	-	ND	-		
50594-66-6	Acifluorfen [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	11 3, 1	D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	- ND	ND	-		
107-02-8	Acrolein [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	1, 3, 3	D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ng Mean for	· Monitoring W	'ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	r Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND			
107-13-1	Acrylonitrile [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
050445 44.0	ADONA I III	C-Zone	ND	ND	ND		ND	ND	ND	ND	ND	ND	NB	ND
958445-44-8	ADONA [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
15072 60 0	Alaskian (v. v. // 1	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
15972-60-8	Alachlor [μg/L]	D-Zone	ND	ND	ND	ND	ND	טא	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	ı	ND			
142363-53-9	Alachlor ESA [µg/L]	C-Zone	ND	0.021	ND	ND	0.00091	ND	ND	0.021	ND	ND	0.00091	ND
142303-33-9	Alactiloi ESA [µg/L]	D-Zone	NA	NA	NA	IND	0.00091	ND	NA	NA	NA	ND	0.00091	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
171262-17-2	Alachlor OA [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
171202 17 2	Aluchiof OA [µg/L]	D-Zone	NA	NA	NA	, ND	IND	ND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
116-06-3	Aldicarb [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
110 00 3	, παιεαί δ [μg/ -]	D-Zone	ND	ND	ND			115	ND	ND	ND	112		113
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	-			ND	-	ND			
1646-88-4	Aldicarb Sulfone [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	i i i i i i i i i i i i i i i i i i i	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	-			ND	-	ND			
1646-87-3	Aldicarb sulfoxide [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND	<u> </u>			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND ND	- ND	ND	-			ND ND	- ND	ND ND			
309-00-2	Aldrin [µg/L]	C-Zone D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
		Sunnyside	ND ND	ND ND	ND ND	-			ND ND	ND ND	ND ND			
		B-Zone	400000	ND	430000				370000	- ND	430000			
		C-Zone	380000	250000	360000	1			330000	230000	340000			
ALK	Alkalinity (as CaCO3) [μg/L]	D-Zone	250000	250000	250000	310000	230000	240000	250000	250000	250000	290000	230000	240000
		Sunnyside	230000	230000	230000	1			230000	230000	230000			
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND ND	63000	ND	1			ND ND	25000	ND ND			
ALKC-1	Alkalinity, Carbonate as CaCO3 [µg/L]	D-Zone	ND ND	ND	ND	ND	2700	ND	ND	ND	ND ND	ND	1100	ND
		Sunnyside	ND	ND	ND	1			ND	ND	ND			







				Us	ing UCL95 f	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	r Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND	-		
ALKH-1	Alkalinity, Hydroxide as CaCO3 [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND	1		
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND			
107-05-1	Allyl Chloride [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
240.04.6	11 2005 43	C-Zone	ND	ND	ND	1	ND		ND	ND	ND	1		ND
319-84-6	alpha-BHC [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
5402.74.0		C-Zone	ND	ND	ND	l NB	ND	ND	ND	ND	ND		NB	ND
5103-71-9	alpha-Chlordane [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	17.79	-	2800				10.17	-	2800			
7420 00 5	Al aria as Tatal Car (1)	C-Zone	60	120	24	1.,	F 2	F1 2	34.43	33.55	15.7	0.0	4.5	50.0
7429-90-5	Aluminum, Total [µg/L]	D-Zone	NA	NA	NA	15.4	5.2	51.3	NA	NA	NA	8.8	1.5	50.9
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	33	-	ND				33	-	ND			
7664-41-7	Ammonio fue // 1	C-Zone	150	1200	ND	25	F2	ND	120	680	ND	29	29	ND
7004-41-7	Ammonia [µg/L]	D-Zone	NA	NA	NA	35	52	ND	NA	NA	NA	29	29	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
26787-78-0	Amoxicillin [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
20707-70-0	Απολιείιιπ [μg/τ]	D-Zone	NA	NA	NA	IND	ND	ND	NA	NA	NA	IND	IND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
62-53-3	Aniline [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
0L 33 3	, ιιπιτε [μg/ -]	D-Zone	ND	ND	ND		ND ND	110	ND	ND	ND		110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	_	ND			
120-12-7	Anthracene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	, w.t dee [p.g/ =]	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.11	-	0.13				0.094	-	0.13			
7440-36-0	Antimony [µg/L]	C-Zone	0.091	0.095	0.12	0.15	0.21	0.21	0.085	0.074	0.11	0.15	0.21	0.21
	, ii J. 1	D-Zone	0.15	0.15	0.15	1			0.15	0.15	0.15			
		Sunnyside	0.22	0.22	0.22				0.22	0.22	0.22			
		B-Zone	ND	-	ND				ND	-	ND			
12674-11-2	Aroclor 1016 [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	i 3. 3	D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND			]	ND	ND	ND			







				Us	ing UCL95 f	or Monitoring V	Vell Data*			Usi	ng Mean for	· Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
11104-28-2	Aroclor 1221 [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
11104 20 2	Αιθείδι 1221 [μg/ Ε]	D-Zone	ND	ND	ND	l ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	  -			ND	-	ND			
11141-16-5	Aroclor 1232 [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	1 LF 3/ 1	D-Zone	ND	ND	ND	-			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND	-			ND	-	ND			
53469-21-9	Aroclor 1242 [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND ND	ND	ND ND	-			ND ND	ND ND	ND ND			
		Sunnyside	ND ND	ND	+				ND ND					
		B-Zone C-Zone	ND ND	- ND	ND ND	1			ND ND	- ND	ND ND			
12672-29-6	Aroclor 1248 [µg/L]	D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND	ND	ND	ND
		Sunnyside	ND ND	ND ND	ND ND	-			ND ND	ND ND	ND			
		B-Zone	ND ND	-	ND ND				ND ND	-	ND			
		C-Zone	ND ND	ND	ND ND	-			ND ND	ND	ND			
11097-69-1	Aroclor 1254 [µg/L]	D-Zone	ND ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND	1			ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND	-			ND	ND	ND			
11096-82-5	Aroclor 1260 [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
22544.54.4		C-Zone	0.97	0.096	ND	1	4.2		0.59	0.096	ND	1.0	4.0	
22541-54-4	Arsenic III [µg/L]	D-Zone	0.39	0.39	0.39	1.9	4.3	4.1	0.23	0.23	0.23	1.8	4.3	4.1
		Sunnyside	4.7	4.7	4.7				4.7	4.7	4.7			
		B-Zone	1.078	-	0.84				0.801	-	0.84			
17428-41-0	Arconic V [ug/l]	C-Zone	1.093	0.73	0.33	0.51	0.032	0.034	0.874	0.46	0.33	0.39	0.02	0.034
17420-41-0	Arsenic V [µg/L]	D-Zone	NA	NA	NA	0.51	0.032	0.034	NA	NA	NA	0.39	0.02	0.034
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.978	-	2.5				0.79	-	2.5			
7440-38-2	Arsenic, Total [μg/L]	C-Zone	1.7	38	1.9	0.58	1.64	0.16	1.393	7.572	1.42	0.47	0.33	0.13
7440 30 2	Alsellie, Total [µg/ L]	D-Zone	NA	NA	NA	0.50	1.04	0.10	NA	NA	NA	0.47	0.55	0.15
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND	_			ND	-	ND			
1332-21-4	Asbestos [MFL]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	. 5555155 [ 2]	D-Zone	ND	ND	ND	1			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.0066	-	ND	-			0.0021	-	ND			
29122-68-7	Atenolol [μg/L]	C-Zone	ND	0.00032	ND	0.002	1.4E-05	ND	ND NA	0.00032	ND	0.0006	1.4E-05	ND
	3	D-Zone Sunnyside	NA NA	NA NA	NA NA	-			NA NA	NA NA	NA NA			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concenti	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production	Concent	ation within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
124522.00.5	Atomorphotics [con/1]	C-Zone	0.00076	ND	ND	0.0001	ND	ND	0.00076	ND	ND	0.0001	ND	ND
134523-00-5	Atorvastatin [μg/L]	D-Zone	NA	NA	NA	0.0001	ND	ND	NA	NA	NA	0.0001	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
1912-24-9	Atrazine [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1912-24-9	Attazine [µg/L]	D-Zone	ND	ND	ND	IND	ND	ND	ND	ND	ND	I ND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.0063	-	ND				0.0063	-	ND			
83905-01-5	Azithromycin [µg/L]	C-Zone	0.015	ND	ND	0.0044	ND	ND	0.012	ND	ND	0.0039	ND	ND
03303 01 3	Azitiiioiiiyeiii [μg/ L]	D-Zone	NA	NA	NA	0.0044	ND	IND	NA	NA	NA	0.0033	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
103-33-3	Azobenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
103 33 3	Α20001126110 [μg/ 2]	D-Zone	NA	NA	NA		ND	IND	NA	NA	NA		IND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	38.98	-	35	-			33.2	-	35			
7440-39-3	Barium, Dissolved [μg/L]	C-Zone	48	19	57	19.8	0.82	4	36	13.13	43.5	16	0.57	3.2
7 110 33 3	Σαπαπη, Σισσόνοα [μg/.2]	D-Zone	NA	NA	NA		0.02		NA	NA	NA		0.57	3.2
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	37.85	-	52	-			34.29	-	52	_		
7440-39-3	Barium, Total [µg/L]	C-Zone	54	29	58	20.4	1.26	4.3	45.8	16.97	45.5	18	0.73	3.6
	- aa, . a aa. [[a-g], -]	D-Zone	NA	NA	NA				NA	NA	NA	_	55	
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND	-			ND	-	ND	1		
25057-89-0	Bentazon [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	11 37 1	D-Zone	ND	ND	ND	-			ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	-			ND	-	ND	-		
100-52-7	Benzaldehyde [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	,	D-Zone	NA	NA	NA	-			NA	NA	NA	-		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND 0.23	- 1.C	ND	-			ND 0.22	1.2	ND	-		
71-43-2	Benzene [µg/L]	C-Zone		1.6	ND	0.039	0.068	ND	0.23 ND	1.2 ND	ND	0.039	0.051	ND
		D-Zone	ND	ND	ND	-			ND ND		ND	-		
		Sunnyside B-Zone	ND ND	ND -	ND ND				ND ND	ND	ND ND			
			UD ND	- ND	ND ND	1			ND ND	- ND	ND ND	+		
92-87-5	Benzidine [µg/L]	C-Zone D-Zone	UD ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
			ND ND	ND ND	ND ND	-			ND ND	ND ND	ND ND	+		
		Sunnyside B-Zone	ND ND	- -	ND ND				ND ND	- -	ND ND			
		C-Zone	ND ND	- ND	ND ND	1			ND ND	- ND	ND ND	1		
56-55-3	Benzo[a]anthracene [µg/L]	D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
			ND ND	ND ND	ND ND	-			ND ND	ND ND	ND ND	+		
		Sunnyside	טא	עא	טא				שו	שוו	עא			







				Us	ing UCL95 fo	or Monitoring \	Well Data*			Us	ing Mean for	Monitoring W	'ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concent	ation within Zone	Capture	Estimated C	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
50.33.0	5	C-Zone	ND	ND	ND	1			ND	ND	ND		ND	
50-32-8	Benzo[a]pyrene [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	1	ND			
205-99-2	Benzo[b]fluoranthene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
205-99-2	венго[р]ниотантнене [µg/L]	D-Zone	ND	ND	ND	IND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
191-24-2	Benzo[g,h,i]perylene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
191-24-2	benzo[g,n,ŋperyiene [µg/L]	D-Zone	ND	ND	ND	IND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
207-08-9	Benzo[k]fluoranthene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
201 00 3	benzo[κ]ndoranthene [μg/ L]	D-Zone	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
65-85-0	Benzoic Acid [μg/L]	C-Zone	ND	6.7	ND	ND	0.29	ND	ND	6.6	ND	ND	0.28	ND
03 03 0	Delizate Acid [µg/ L]	D-Zone	ND	ND	ND		0.23	IND	ND	ND	ND	ND	0.20	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	NA	-	NA				NA	-	NA			
95-16-9	Benzothiazole [µg/L]	C-Zone	NA	8.5	NA	NA	0.37	NA	NA	6.3	NA	NA	0.27	NA
	Joi. 1201. 102-0-1 [p.g/ -]	D-Zone	NA	NA	NA		0.0.		NA	NA	NA		0.2.	
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND	_			ND	-	ND			
100-51-6	Benzyl Alcohol [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	, , , , , , , , , , , , , , , , , , ,	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	1.1	-	6.3	1			0.69	-	6.3			
85-68-7	Benzyl butyl phthalate [µg/L]	C-Zone	0.69	0.44	0.53	0.51	0.17	0.29	0.54	0.3	0.53	0.36	0.17	0.29
	, , ,	D-Zone	ND 0.17	ND	ND 0.17	1			ND	ND	ND			
		Sunnyside	0.17	0.17	0.17				0.17	0.17	0.17			
		B-Zone	ND	- ND	ND	-			ND	- ND	ND			
7440-41-7	Beryllium [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	· -	D-Zone	ND	ND	ND	-			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone C-Zone	ND ND	- ND	ND ND	1			ND ND	- ND	ND ND			
319-85-7	beta-BHC [μg/L]	D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND	ND	ND	ND
			ND ND		ND ND	+			ND ND	ND ND	ND			
		Sunnyside B-Zone	510000	ND -	520000				450000	- -	520000			
		C-Zone	480000	280000	430000	+			410000	250000	410000			
71-52-3	Bicarbonate Alkalinity as HCO3 [µg/L]	D-Zone	310000	310000	310000	390000	280000	290000	310000	310000	310000	360000	280000	290000
			280000	280000	280000	+			280000	280000				
		Sunnyside	200000	200000	200000				200000	200000	280000			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
21.2511			Concent	ration within	Capture	Estimated Co	oncentrations fo	or Production	Concent	ration within			oncentrations f	or Production
CASRN	Constituent [Units]	Layer		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
BOD	Biochemical Oxygen Demand [µg/L]	C-Zone	ND	6900	ND	ND	300	ND	ND	5300	ND	ND	230	ND
ВОО	Biochemical Oxygen Demand [μg/L]	D-Zone	NA	NA	NA	ND	300	IND	NA	NA	NA	IND	230	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
111-91-1	bis(2-chloroethoxy)methane [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
111-31-1	bis(2-chioroethoxy)methane [µg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	IND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
111-44-4	bis(2-chloroethyl)ether [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
111-44-4	bis(2-ciliordethyr)ether [µg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	IND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
108-60-1	bis(2-chloroisopropyl)ether [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
100-00-1	bis(z-cilioroisopropyr)etrier [µg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	IND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
103-23-1	Bis(2-ethylhexyl)adipate [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
103-23-1	bis(2-etilyinexyi)adipate [µg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	IND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
117-81-7	bis(2-ethylhexyl)phthalate [µg/L]	C-Zone	ND	0.91	ND	0.38	0.7	0.63	ND	0.73	ND	0.35	0.68	0.62
117-01-7	bis(2-etifyfflexyf)pfftflafate [µg/L]	D-Zone	0.78	0.78	0.78	0.36	0.7	0.03	0.62	0.62	0.62	0.55	0.00	0.02
		Sunnyside	0.68	0.68	0.68				0.68	0.68	0.68			
		B-Zone	0.0052	-	0.00053				0.0032	-	0.00053			
80-05-7	Bisphenol A [µg/L]	C-Zone	0.0082	0.13	0.0023	0.0029	0.0056	0.0001	0.0034	0.066	0.0022	0.0015	0.0029	0.0001
00 03 7	ызрненог д [µg/ с]	D-Zone	NA	NA	NA	0.0023	0.0030	0.0001	NA	NA	NA	0.0013	0.0023	0.0001
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	143.6	-	180				135	-	180			
7440-42-8	Boron, Total [μg/L]	C-Zone	119.6	600	140	63.3	26	11.4	110	298	120	59.1	12.9	10.3
7440 42 0	ουτοτί, τοται [μg/ <u>Ε</u> ]	D-Zone	ND	ND	ND	03.3	20	11	ND	ND	ND	33.1	12.3	10.5
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
314-40-9	Bromacil [µg/L]	C-Zone	0.37	ND	ND	0.062	ND	ND	0.37	ND	ND	0.062	ND	ND
314 40 3	Dromacii [μg/ L]	D-Zone	ND	ND	ND	0.002	145	110	ND	ND	ND	0.002	110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	28	-	ND				13	-	ND			
15541-45-4	Bromate [µg/L]	C-Zone	7.7	160	10	10	7.1	0.76	7.7	52	10	5.7	2.4	0.76
13311 43 4	Σ.σαιο [μg/ L]	D-Zone	3	3	3			0.70	3	3	3	]		3.70
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	790	-	440				670	-	440			
24959-67-9	Bromide [µg/L]	C-Zone	560	5700	400	330	250	31	400	2200	350	270	97	28
24333 01-3	brofflide [μg/ L]	D-Zone	NA	NA	NA	330	250		NA	NA	NA			20
		Sunnyside	NA	NA	NA				NA	NA	NA			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	· Monitoring W	ell Data**	
CASRN	Constituent [Unite]	Layer	Concent	ration within	Capture	Estimated C	oncentrations fo	or Production	Concent	ration within	Capture	Estimated Co	oncentrations fo	or Production
CASKIN	Constituent [Units]	Layer		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
108-86-1	Bromobenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
100-00-1	втотпорендене [µg/L]	D-Zone	ND	ND	ND	ND	IND	IND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
5589-96-8	Bromochloroacetic acid [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
3309-90-0	βιοιποεποιοάcetic acid [μg/L]	D-Zone	ND	ND	ND	ND	IND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
74-97-5	Bromochloromethane [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
14-91-5	втотпостноготпентале [µg/L]	D-Zone	ND	ND	ND	ND	IND	IND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	NA	-	NA				NA	-	NA			
71133-14-7	Bromodichloroacetic acid [µg/L]	C-Zone	NA	ND	NA	ND	ND	ND	NA	ND	NA	ND	ND	ND
/1135-14-/	bromodichioroacetic acid [µg/L]	D-Zone	ND	ND	ND	ND	IND	IND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.26	-	ND				0.24	-	ND			
75-27-4	Bromodichloromethane [µg/L]	C-Zone	ND	ND	ND	0.077	ND	ND	ND	ND	ND	0.072	ND	ND
15-21-4	Biomodichioromethane [μg/L]	D-Zone	ND	ND	ND	0.077	IND	IND	ND	ND	ND	0.072	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
75-25-2	Promoform [ug/l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
15-25-2	Bromoform [µg/L]	D-Zone	ND	ND	ND	ND	IND	IND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
74-83-9	Promomethano [ug/l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
74-05-9	Bromomethane [μg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	שוו	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
23184-66-9	Butachlor [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
23104-00-9	butacinoi [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	1.6	_	ND				1.6	-	ND			
123-72-8	Butanal [μg/L]	C-Zone	2	ND	ND	0.82	ND	ND	2	ND	ND	0.82	ND	ND
123-12-0	butanai [µg/L]	D-Zone	NA	NA	NA	0.62	IND	ND	NA	NA	NA	0.02	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.094	-	ND				0.058	-	ND			
7440-43-9	Cadmium [µg/L]	C-Zone	0.073	ND	ND	0.054	0.004	0.0041	0.073	ND	ND	0.041	0.0035	0.0035
1770-45-3	Caumum [μg/L]	D-Zone	0.072	0.072	0.072	0.054	0.004	0.0041	0.063	0.063	0.063	0.041	0.0033	0.0033
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.031	-	0.0013				0.016	-	0.0013			
58-08-2	Caffeine [µg/L]	C-Zone	0.0096	0.06	0.0038	0.011	0.0026	0.0003	0.0039	0.02	0.0027	0.0056	0.00084	0.0002
30-00-2	Carrelle [µg/L]	D-Zone	ND	ND	ND	0.011	0.0020	0.0003	ND	ND	ND	0.0030	0.0004	0.0002
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ng Mean for	· Monitoring W	ell Data**	
CACDN	Compliance Illustra		Concenti	ation within	Capture	Estimated C	oncentrations fo	or Production	Concentr	ation within	Capture	Estimated C	oncentrations fo	or Production
CASRN	Constituent [Units]	Layer		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	180000	-	140000				150000	-	140000			
7440-70-2	Calcium [µg/L]	C-Zone	150000	100000	130000	120000	73000	76000	120000	63000	130000	110000	71000	76000
1440 70 2	Calciam [µg/L]	D-Zone	95000	95000	95000	120000	73000	70000	94000	94000	94000	110000	7 1000	70000
		Sunnyside	70000	70000	70000				70000	70000	70000			
		B-Zone	ND	-	ND	-			ND	-	ND			
133-06-2	Captan [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
.55 55 2		D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	-			ND	-	ND			
298-46-4	Carbamazepine [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	22 22 2 2 2 E 3, 1	D-Zone	NA	NA	NA	-			NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND	-			ND	-	ND			
63-25-2	Carbaryl [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	, 3	D-Zone	ND	ND	ND	-			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	-			ND	- ND	ND			
86-74-8	Carbazole [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	-	D-Zone	ND	ND	ND	-			ND	ND	ND			
		Sunnyside	ND ND	ND	ND				ND ND	ND	ND ND			
		B-Zone C-Zone	ND ND	- ND	ND ND	1			ND	- ND	ND ND			
1563-66-2	Carbofuran [µg/L]	D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND	ND ND	ND ND	ND	ND	ND
		Sunnyside	ND ND	ND	ND ND	-			ND	ND ND	ND ND			
		B-Zone	ND ND	-	ND ND				ND	-	ND			
		C-Zone	ND	0.26	ND				ND	0.26	ND			
75-15-0	Carbon Disulfide [µg/L]	D-Zone	ND	ND	ND	ND	0.011	ND	ND	ND	ND	ND	0.011	ND
		Sunnyside	ND	ND	ND	1			ND	ND	ND			
		B-Zone	0.63	-	ND				0.512	-	ND			
		C-Zone	0.675	0.62	0.12	-			0.571	0.517	0.12			
56-23-5	Carbon Tetrachloride [µg/L]	D-Zone	0.3	0.3	0.3	0.36	0.044	0.024	0.3	0.3	0.3	0.3	0.039	0.024
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
100.00		C-Zone	ND	ND	ND	1			ND	ND	ND			
133-90-4	Chloramben [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	131.8	-	160				98.67	-	160			
2146052	Chloreta Francis	C-Zone	86.96	33	140	[ [ [ ]	1.4	11 1	86	23.5	92.5	441	1	0.3
2146053	Chlorate [μg/L]	D-Zone	NA	NA	NA	54.3	1.4	11.1	NA	NA	NA	44.1	'	8.3
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
57-74-9	Chlordane [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
31-14-9	Chlordane [µg/L]	D-Zone	ND	ND	ND	שמו	שויו	שוו	ND	ND	ND	אט	שוו	שוו
		Sunnyside	ND	ND	ND				ND	ND	ND			<u> </u>







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ng Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concent	ation within Zone	Capture	Estimated Co	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	130000	-	86000				120000	1	86000			
16007.00.6		C-Zone	110000	600000	93000	02000	0.000	65000	100000	280000	82000	0.000	72000	64000
16887-00-6	Chloride [µg/L]	D-Zone	74000	74000	74000	93000	86000	65000	70000	70000	70000	86000	72000	64000
		Sunnyside	62000	62000	62000				62000	62000	62000			
		B-Zone	ND	-	ND				ND	-	ND			
76-13-1	Chlorinated Fluorocarbon (Freon 113) [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
70-13-1	Chiomiated Fidorocarbon (Freon 113) [µg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	IND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	6	-	56				4.3	-	56			
7782-50-5RF	Chlorine Residual, Free [µg/L]	C-Zone	15	13	80	4.3	0.56	5.7	15	11	48	3.7	0.48	3.8
1102 30 SKI	Chlorine Residual, Free [µg/L]	D-Zone	NA	NA	NA	4.5	0.50	5.7	NA	NA	NA	3.7	0.40	3.0
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	20	-	160				14	-	160			
TOTAL-	Chlorine Residual, Total [µg/L]	C-Zone	29	22	100	11	0.95	8.7	24	20	61	8.3	0.87	6.4
CHLORINE	emorme Nesidadi, Fotal [µg, 2]	D-Zone	NA	NA	NA		0.55	0.7	NA	NA	NA		0.07	0.4
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND	<u> </u>  -		
14998-27-7	Chlorite [µg/L]	C-Zone	ND	16	ND	1.6	3.5	2.7	ND	14	ND	1.6	3.4	2.7
		D-Zone	3.3	3.3	3.3				3.1	3.1	3.1			
		Sunnyside	2.9	2.9	2.9				2.9	2.9	2.9			
		B-Zone	ND	-	ND				ND	-	ND	-		
79-11-8	Chloroacetic Acid [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	J	D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- 100	ND				ND	-	ND	-		
108-90-7	Chlorobenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND ND	ND	ND				ND ND	ND	ND ND	-		
		Sunnyside	NA	ND	ND NA				NA NA	ND	NA NA			
		B-Zone C-Zone	NA NA	- ND	NA NA				NA NA	- ND	NA NA	1		
5278-95-5	Chlorodibromoacetic acid [µg/L]	D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND	ND	ND ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND	-		
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND	1		
75-00-3	Chloroethane [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND	1		
		B-Zone	1.4	-	19				1.3	-	15			
		C-Zone	2.7	2.5	1.1				2.2	2.1	0.86	1		
67-66-3	Chloroform [µg/L]	D-Zone	1.3	1.3	1.3	1.1	0.18	0.47	1	1	1	0.95	0.15	0.37
		Sunnyside	ND	ND	ND				ND	ND	ND	=		
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND	1		
74-87-3	Chloromethane [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND	1		







						or Monitoring V						r Monitoring W		
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concent	ation within Zone	Capture	Estimated C	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
1897-45-6	Chlorothalonil [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1097-43-0	Chlorothalothi [µg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND		ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND	_		
101-21-3	Chlorpropham [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	F - F	D-Zone	ND	ND	ND				ND	ND	ND	1		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.68	- 0.016	5.9				0.26	- 0.01.1	5.9	-		
18540-29-9	Chromium, Hexavalent [µg/L]	C-Zone	1.5	0.016	4.5	0.46	0.017	0.39	0.84	0.014	2.9	0.23	0.017	0.29
		D-Zone	0.0092	0.0092	0.0092				0.0092	0.0092	0.0092	-		
		Sunnyside	0.018 9.905	0.018	0.018 17				0.018	0.018	0.018 17			
		B-Zone C-Zone	290	1.3	240				2.318 145.9	0.873	127.5	+		
7440-47-3	Chromium, Total [μg/L]	D-Zone	NA	NA	NA	51.5	0.056	14.4	NA	NA	NA	25.1	0.038	7.8
		Sunnyside	NA NA	NA	NA NA				NA	NA NA	NA NA	1		
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND ND	ND	ND				ND	ND	ND	1		
218-01-9	Chrysene [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.0062	-	0.0045				0.0042	-	0.0045			
		C-Zone	0.014	0.0027	0.013				0.0057	0.0027	0.0076	1		
85721-33-1	Ciprofloxacin [µg/L]	D-Zone	NA	NA	NA	0.0042	0.00012	0.0008	NA	NA	NA	0.0022	0.00012	0.0005
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.26	-	ND				0.26	-	ND			
156 50 2		C-Zone	0.852	49.37	0.8	0.22	2.1	0.047	0.736	30.44	0.613	0.2	1.2	0.026
156-59-2	cis-1,2-Dichloroethene (cis-1,2-DCE) [μg/L]	D-Zone	ND	ND	ND	0.22	2.1	0.047	ND	ND	ND	0.2	1.3	0.036
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
10061-01-5	cis-1,3-Dichloropropene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
10001-01-3	cis-1,3-Dictiloroproperie [µg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	IND	IND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
5103-73-1	cis-Nonachlor [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
3103 73 1	εισ (νοτιαετιίοι [μg/ <u>Ε</u> ]	D-Zone	ND	ND	ND	110	110	110	ND	ND	ND		145	ND ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	7.6	-	0.014				1.8	-	0.014	_		
7440-48-4	Cobalt [µg/L]	C-Zone	2.6	0.093	1.3	2.7	0.004	0.077	1.3	0.093	0.98	0.76	0.004	0.057
	1	D-Zone	NA	NA	NA				NA	NA	NA	1		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	- ND	ND				ND	-	ND	-		
CLR	Color [Units]	C-Zone	ND	ND ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Us	ing UCL95 fo	or Monitoring V	/ell Data*	Using Mean for Monitoring Well Data**						
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production	Concenti	ation within Zone	Capture	Estimated Concentrations for Productions Wells		
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
CLD	Calan II Inital	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
CLR	Color [Units]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.96	-	0.49				0.47	-	0.49			
7440-50-8	Copper [ug/l]	C-Zone	2.4	0.23	1.6	0.84	0.22	0.3	1	0.23	1.1	0.46	0.21	0.27
7440-30-6	Copper [µg/L]	D-Zone	0.45	0.45	0.45	0.04	0.22	0.3	0.42	0.42	0.42	0.40	0.21	0.27
		Sunnyside	0.2	0.2	0.2				0.2	0.2	0.2			
		B-Zone	0.0028	-	0.00065				0.0013	-	0.00065			
486-56-6	Cotinine [µg/L]	C-Zone	0.004	0.027	0.0026	0.0015	0.0012	0.0002	0.0029	0.015	0.0018	0.0009	0.00063	0.0001
400-30-0	Courinie [µg/L]	D-Zone	NA	NA	NA	0.0013	0.0012	0.0002	NA	NA	NA	0.0003	0.00003	0.0001
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
123-73-9	Crotonaldehyde [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
125 75 5	Crotonaldenyde [µg/L]	D-Zone	NA	NA	NA	ND	ND	IND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
21725-46-2	Cyanazine [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
21723 40 2	Cydnaznie [µg/ L]	D-Zone	ND	ND	ND	110	110	145	ND	ND	ND	110	110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
57-12-5	Cyanide [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
3, 123	εγαιπαε [μg/ -]	D-Zone	ND	ND	ND	145	115		ND	ND	ND	145	1,15	115
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	NA	-	NA				NA	-	NA			
1122-82-3	Cyclohexane, isothiocyanato- [µg/L]	C-Zone	NA	1	NA	NA	0.043	NA	NA	1	NA	NA	0.043	NA
	cy c.e	D-Zone	NA	NA	NA		0.0.0		NA	NA	NA		0.0.0	
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
108-94-1	Cyclohexanone [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	-y	D-Zone	NA	NA	NA				NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
75-99-0	Dalapon [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	,3	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	- ND	ND			
1861-32-1	DCPA [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	- ND	ND			
112-31-2	Decanal [μg/L]	C-Zone	ND	ND NA	ND	ND	ND	ND	ND NA	ND	ND	ND	ND	ND
		D-Zone	NA	NA NA	NA				NA NA	NA	NA NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			







				Us	ing UCL95 f	or Monitoring V	Vell Data*			Usi	ing Mean foi	r Monitoring W	/ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within	Capture	Estimated C	oncentrations fo	r Production	Concentr	ation within	Capture	Estimated C	oncentrations fo	r Production
G.151				Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	0.0031	-	0.0016	-			0.002	-	0.0016			
134-62-3	DEET [μg/L]	C-Zone	0.0011	0.0036	0.0027	0.0011	0.00016	0.0002	0.0006	0.0033	0.0024	0.0007	0.00014	0.0002
		D-Zone	NA	NA NA	NA	-			NA NA	NA NA	NA NA			
		Sunnyside B-Zone	NA ND	NA -	NA ND				NA ND	NA -	NA ND			
		C-Zone	ND ND	ND	ND				ND	ND	ND			
319-86-8	delta-BHC [µg/L]	D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND	ND ND	ND	ND	ND	ND
		Sunnyside	ND ND	ND	ND ND	-			ND	ND ND	ND			
		B-Zone	0.00021	-	ND				0.00021	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND			
439-14-5	Diazepam [μg/L]	D-Zone	NA	NA	NA	6E-05	ND	ND	NA	NA	NA	6E-05	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
222 44 5	D: :	C-Zone	ND	0.055	ND		0.0004	ND	ND	0.055	ND	ND	0.0004	NB
333-41-5	Diazinon [μg/L]	D-Zone	ND	ND	ND	ND	0.0024	ND	ND	ND	ND	ND	0.0024	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
53-70-3	Dibenzo[a,h]anthracene [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
33 70 3	Dibenzo[α,njantinαcene [μg/ ε]	D-Zone	ND	ND	ND		ND	ND	ND	ND	ND	ND	IVD	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	-			ND	-	ND			
132-64-9	Dibenzofuran [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	1, 3, 1	D-Zone	ND	ND	ND	-			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND ND	- ND	ND ND	-			ND ND	- ND	ND ND			
631-64-1	Dibromoacetic acid [µg/L]	C-Zone D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
		Sunnyside	ND ND	ND ND	ND ND	-			ND	ND ND	ND ND			
		B-Zone	ND ND	-	ND				ND	-	ND ND			
		C-Zone	ND	70	ND	-			ND	0.86	ND			
124-48-1	Dibromochloromethane [µg/L]	D-Zone	ND	ND	ND	ND	3	ND	ND	ND	ND	ND	0.037	ND
		Sunnyside	ND	ND	ND	-			ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
	<b>-</b> 0	C-Zone	ND	ND	ND	1			ND	ND	ND			
74-95-3	Dibromomethane [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
1918-00-9	Dicamba [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1910-00-9	ысаныа [ру/г]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	ND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	4.5	-	ND	_			3	-	ND			
3400-09-7	Dichloramine [µg/L]	C-Zone	6	4	15	2.4	0.17	0.88	5	4	8.5	1.7	0.17	0.5
	2.6 [pg/ 2]	D-Zone	NA	NA	NA			2.00	NA	NA	NA	,		3.3
		Sunnyside	NA	NA	NA				NA	NA	NA			







				Us	ing UCL95 f	or Monitoring \	Well Data*		Using Mean for Monitoring Well Data**							
CASRN	Constituent [Units]	Layer	Concenti	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concentr	ation within Zone	Capture	Estimated Concentrations for Production Wells				
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9		
		B-Zone	ND	-	ND				ND	-	ND					
		C-Zone	ND	ND	ND	1			ND	ND	ND					
79-43-6	Dichloroacetic Acid [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	=	ND					
75 71 0	D'able es d'O e constitue es for (1)	C-Zone	ND	ND	ND	1	ND	ND	ND	ND	ND	ND	ND	ND		
75-71-8	Dichlorodifluoromethane [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
120.26.5	D' 11	C-Zone	ND	ND	ND	1	NB	ND	ND	ND	ND	ND	NB	ND		
120-36-5	Dichloroprop [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	0.00043	-	ND				0.00043	-	ND					
15207.06.5	D'alafa a a f	C-Zone	0.00073	ND	ND	0.0003	ND	ND	0.00073	ND	ND	0.0003	ND	ND		
15307-86-5	Diclofenac [μg/L]	D-Zone	NA	NA	NA	0.0003	ND	ND	NA	NA	NA	0.0003	ND	ND		
		Sunnyside	NA	NA	NA				NA	NA	NA					
		B-Zone	ND	-	ND				ND	-	ND					
60 57 1	District of the	C-Zone	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND		
60-57-1	Dieldrin [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	0.093	-	ND				0.093	-	ND					
04.66.2	D'ada Lababadara La #1	C-Zone	0.1	0.63	0.084	0.14	0.10	0.16	0.1	0.31	0.084	0.12	0.16	0.15		
84-66-2	Diethyl phthalate [μg/L]	D-Zone	0.17	0.17	0.17	0.14	0.19	0.16	0.13	0.13	0.13	0.12	0.16	0.15		
		Sunnyside	0.17	0.17	0.17				0.16	0.16	0.16					
		B-Zone	ND	-	ND				ND	-	ND					
FC F2 1	Diate latille actual from /11	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
56-53-1	Diethylstilbestrol [µg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA	ND	ND	ND		
		Sunnyside	NA	NA	NA				NA	NA	NA					
		B-Zone	ND	-	ND				ND	-	ND					
100 20 2	Di isanganul athar (um/l)	C-Zone	ND	1.2	ND	ND	0.052	ND	ND	1.1	ND	ND	0.05	ND		
108-20-3	Di-isopropyl ether [μg/L]	D-Zone	ND	ND	ND	ND	0.053	ND	ND	ND	ND	ND	0.05	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
60-51-5	Dimenth acts [ug/l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
00-51-5	Dimethoate [μg/L]	D-Zone	ND	ND	ND	IND	IND	ND	ND	ND	ND	ND	שוו	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
131-11-3	Dimethyl phthalate [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
151-11-5	Difficulty prictialate [µg/L]	D-Zone	ND	ND	ND	IND	IND	ND	ND	ND	ND	ND	IND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	0.28	-	2.7				0.28	-	2.7					
84-74-2	Di n hutul phthalata (ua/li)	C-Zone	0.25	0.26	0.23	0.26	0.25	0.29	0.23	0.26	0.21	0.25	0.25	0.29		
04-74-2	Di-n-butyl-phthalate [µg/L]	D-Zone	0.26	0.26	0.26	0.26	0.25	0.29	0.24	0.24	0.24	0.25	0.25	0.29		
		Sunnyside	0.25	0.25	0.25				0.25	0.25	0.25					







				Us	ing UCL95 fo	or Monitoring V	Vell Data*		Using Mean for Monitoring Well Data**							
CASRN	Constituent [Units]	Layer	Concent	ration within Zone			oncentrations fo Wells	or Production	Concent	ration within Zone			oncentrations fo Wells	or Production		
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9		
		B-Zone	ND	-	ND				ND	-	ND					
117.04.0	D' e est l'abilitate ( e //)	C-Zone	ND	ND	ND	, ND	ND	ND	ND	ND	ND	l NB	ND	ND		
117-84-0	Di-n-octyl phthalate [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
88-85-7	Dinoseb [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
00-03-7	Dinoseb [µg/L]	D-Zone	ND	ND	ND		IND	ND	ND	ND	ND		ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
957-51-7	Diphenamid [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
337-31-7	Бірпепатііц [µg/ с]	D-Zone	ND	ND	ND	I ND	ND	ND	ND	ND	ND		ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
85-00-7	Diquat [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
03 00 7	Diquat [µg/ L]	D-Zone	ND	ND	ND		145	110	ND	ND	ND		110	140		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND	_			ND	-	ND					
298-04-4	Disulfoton [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
	Σισαποτοπ [μg/ []	D-Zone	ND	ND	ND	1			ND	ND	ND	_				
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
330-54-1	Diuron [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
	J	D-Zone	NA	NA	NA				NA	NA	NA					
		Sunnyside	NA NB	NA	NA				NA	NA	NA					
		B-Zone	ND	- ND	ND	_			ND	- ND	ND	_				
959-98-8	Endosulfan I [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		D-Zone	ND	ND ND	ND	_			ND	ND	ND	-				
		Sunnyside	ND ND	ND	ND				ND	ND	ND ND					
		B-Zone C-Zone	ND ND	- ND	ND ND	1			ND ND	- ND	ND	+				
33213-65-9	Endosulfan II [µg/L]	D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND	ND ND	ND ND	ND	ND	ND		
		Sunnyside	ND ND	ND ND	ND ND	-			ND	ND ND	ND ND	-				
		B-Zone	ND ND	-	ND ND				ND	-	ND ND					
		C-Zone	ND ND	ND	ND ND	-			ND	ND	ND ND	-				
1031-07-8	Endosulfan sulfate [µg/L]	D-Zone	ND ND	ND	ND	ND	ND	ND	ND	ND	ND ND	ND	ND	ND		
		Sunnyside	ND ND	ND	ND	-			ND	ND	ND	1				
		B-Zone	ND ND	-	ND				ND	-	ND					
		C-Zone	ND ND	ND	ND				ND	ND	ND					
145-73-3	Endothall [µg/L]	D-Zone	ND ND	ND	ND	ND	ND	ND	ND	ND ND	ND ND	ND	ND	ND		
		Sunnyside	ND ND	ND	ND	1			ND	ND	ND					
		B-Zone	ND ND	-	ND			1	ND	-	ND					
		C-Zone	ND ND	ND	ND	1			ND	ND	ND					
72-20-8	Endrin [µg/L]	D-Zone	ND ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND	1			ND	ND	ND					
		Junnyside	NU	שויו	שוו	I.	<u> </u>	1	שויו	שויו	שוו		<u> </u>			







			Using UCL95 for Monitoring Well Data*							Using Mean for Monitoring Well Data**						
CASRN	Constituent [Units]	Layer	Concent	ration within Zone			oncentrations fo Wells	r Production	Concenti	ation within Zone			oncentrations fo Wells	r Production		
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9		
		B-Zone	ND	-	ND				ND	-	ND					
7424 02 4	5 1: ALL 1 7 (1)	C-Zone	ND	ND	ND		NB	NB	ND	ND	ND	1	ND	ND		
7421-93-4	Endrin Aldehyde [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
53494-70-5	Endrin Ketone [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
33494-70-3	Endrin Retone [µg/L]	D-Zone	ND	ND	ND	ND	IND	IND	ND	ND	ND	IND	ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
481-30-1	Epitestosterone [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
401 30 1	εριτεsτοsτεί οπε [μg/ε]	D-Zone	NA	NA	NA	I ND	ND	ND	NA	NA	NA		ND	ND		
		Sunnyside	NA	NA	NA				NA	NA	NA					
		B-Zone	ND	-	ND				ND	-	ND					
759-94-4	EPTC [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
	Στ. το [μg/ -]	D-Zone	ND	ND	ND		115	1,15	ND	ND	ND					
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
50-27-1	Estriol [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
	LP 3/ 1	D-Zone	NA	NA	NA				NA	NA	NA					
		Sunnyside	NA	NA	NA				NA	NA	NA					
		B-Zone	ND 0.0005	- ND	ND	-			ND 0.0003.6	- ND	ND	<u> </u>				
53-16-7	Estrone [µg/L]	C-Zone	0.0005	ND	0.00077	8E-05	ND	5E-05	0.00036	ND NA	0.00077	6E-05	ND	5E-05		
		D-Zone	NA NA	NA	NA	-			NA	NA NA	NA	<u> </u>				
		Sunnyside	NA	NA	NA				NA	NA	NA					
		B-Zone	ND 0.0016	- ND	ND	-			ND 0.0016	- ND	ND	_				
2991-50-6	EtFOSAA [µg/L]	C-Zone D-Zone	0.0016 ND	ND ND	ND ND	0.0003	ND	ND	0.0016 ND	ND ND	ND ND	0.0003	ND	ND		
		Sunnyside	NA NA	NA NA	NA NA	-			NA NA	NA NA	NA NA	_				
		B-Zone	ND ND	INA -	ND ND				ND ND	- IVA	ND ND					
		C-Zone	ND ND	ND	ND	-			ND	ND	ND					
64-17-5	Ethanol [µg/L]	D-Zone	NA NA	NA NA	NA NA	ND	ND	ND	NA NA	NA NA	NA NA	ND	ND	ND		
		Sunnyside	NA NA	NA NA	NA NA	-			NA NA	NA NA	NA NA	1				
		B-Zone	ND ND	-	ND				ND	-	ND					
		C-Zone	ND ND	ND	ND	-			ND	ND	ND					
563-12-2	Ethion [μg/L]	D-Zone	ND ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND	-			ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
		C-Zone	ND	ND	ND	-			ND	ND	ND					
97-63-2	Ethyl Methacrylate [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND				ND	ND	ND					
		B-Zone	ND	-	ND				ND	-	ND					
		C-Zone	ND	ND	ND	1			ND	ND	ND					
637-92-3	Ethyl tert-butyl ether [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
		Sunnyside	ND	ND	ND	1			ND	ND	ND					







				Us	ing UCL <u>95</u> fo	or Monitoring V	Vell Data*		Using Mean for Monitoring Well Data**						
CASRN	Constituent [Units]	Layer	Concenti	ation within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production	Concent	ation within Zone	Capture	Estimated Co	oncentrations fo Wells	r Production	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	
		B-Zone	ND	-	ND				ND	-	ND				
100 44 4	Ed. II	C-Zone	ND	ND	ND	ND	NB	ND	ND	ND	ND		ND	ND	
100-41-4	Ethylbenzene [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
		Sunnyside	ND	ND	ND				ND	ND	ND				
		B-Zone	ND	-	ND				ND	1	ND				
107-21-1	Ethylana Chysol (us /I.)	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
107-21-1	Ethylene Glycol [µg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA	ND	ND	ND	
		Sunnyside	NA	NA	NA				NA	NA	NA				
		B-Zone	ND	=	ND				ND	-	ND				
206-44-0	Fluoranthene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
200-44-0	Fluorantherie [µg/L]	D-Zone	ND	ND	ND	ND	IND	IND	ND	ND	ND	IND	ND	ND	
		Sunnyside	ND	ND	ND				ND	ND	ND				
		B-Zone	ND	-	ND				ND	-	ND				
86-73-7	Fluorene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
00-73-7	ridorene [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	IND	ND	ND	
		Sunnyside	ND	ND	ND				ND	ND	ND				
		B-Zone	0.307	-	0.34				0.268	-	0.34				
16984-48-8	Fluoride, Total [mg/L]	C-Zone	0.32	0.23	0.31	0.15	0.01	0.024	0.224	0.19	0.3	0.12	0.0082	0.024	
10304 40 0	ridonde, rotai [mg/t]	D-Zone	NA	NA	NA	0.13	0.01	0.024	NA	NA	NA	0.12	0.0002	0.024	
		Sunnyside	NA	NA	NA				NA	NA	NA				
		B-Zone	0.00095	-	ND				0.00061	-	ND	-			
54910-89-3	Fluoxetine [µg/L]	C-Zone	0.00094	ND	0.0002	0.0004	ND	1E-05	0.00039	ND	0.0002	0.0003	ND	1E-05	
34310 03 3	ridoxetine [µg/ L]	D-Zone	NA	NA	NA	0.0004	110	12 03	NA	NA	NA	0.0003	110	12 03	
		Sunnyside	NA	NA	NA				NA	NA	NA				
		B-Zone	ND	_	ND				ND	-	ND	_			
944-22-9	Fonofos [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
311 == 3	. ee.es [µg/ =]	D-Zone	NA	NA	NA				NA	NA	NA		2		
		Sunnyside	NA	NA	NA				NA	NA	NA				
		B-Zone	ND	-	ND				ND	-	ND	-			
50-00-0	Formaldehyde [µg/L]	C-Zone	ND	0.89	ND	ND	0.039	ND	ND	0.89	ND	ND	0.039	ND	
	, i. J. 1	D-Zone	NA	NA	NA				NA	NA	NA	-			
		Sunnyside	NA	NA	NA				NA	NA	NA				
		B-Zone	0.046	-	0.012				0.029	-	0.012	-			
1222-05-5	Galaxolide (HHCB) [μg/L]	C-Zone	0.042	0.024	0.019	0.021	0.001	0.0013	0.025	0.017	0.017	0.013	0.00074	0.0012	
	, , , , , , , , , , , , , , , , , , ,	D-Zone	NA	NA	NA				NA	NA	NA	-			
		Sunnyside	NA	NA	NA				NA	NA	NA				
		B-Zone	ND	- ND	ND				ND	- ND	ND	-			
58-89-9	gamma-BHC [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
		D-Zone	ND	ND	ND				ND	ND	ND	-			
		Sunnyside	ND	ND	ND			1	ND	ND	ND			-	
		B-Zone	ND	- ND	ND				ND	- ND	ND	-			
5566-34-7	gamma-Chlordane [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
		D-Zone	ND	ND	ND				ND	ND	ND	-			
		Sunnyside	ND	ND	ND				ND	ND	ND				







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ing Mean foi	<sup>r</sup> Monitoring W	ell Data**	
CASRN	Constituent [[Inite]	Lavar	Concent	ration within	Capture	Estimated C	oncentrations fo	or Production	Concent	ration within	Capture	Estimated C	oncentrations fo	r Production
CASKN	Constituent [Units]	Layer		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	120	-	100				94	-	100			
8032-32-4	Gasoline Range Organics [µg/L]	C-Zone	860	100	120	180	4.3	8.8	330	100	110	84	4.3	7.9
0032 32 4	Gasoniie Harige Organies [µg/2]	D-Zone	NA	NA	NA	100	4.5	0.0	NA	NA	NA	0-1	4.5	7.5
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.00011	-	ND	-			0.00011	-	ND			
25812-30-0	Gemfibrozil [µg/L]	C-Zone	ND	0.00012	ND	3E-05	5.2E-06	ND	ND	0.00012	ND	3E-05	5.2E-06	ND
	<u></u> 3	D-Zone	NA	NA	NA	-			NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND ND	- ND	ND	-			ND ND	- ND	ND ND			
107-22-2	Glyoxal [µg/L]	C-Zone D-Zone	NA	ND NA	ND NA	ND	ND	ND	NA NA	NA NA	NA NA	ND	ND	ND
		Sunnyside	NA NA	NA NA	NA NA	1			NA NA	NA NA	NA NA			
		B-Zone	ND ND	INA	ND				ND	- INA	ND ND			
		C-Zone	ND ND	2.2	ND	-			ND	2.2	ND			
1071-83-6	Glyphosate [µg/L]	D-Zone	2.1	2.1	2.1	0.38	0.21	0.12	2.1	2.1	2.1	0.38	0.21	0.12
		Sunnyside	ND	ND	ND	-			ND	ND	ND			
		B-Zone	6.663	-	5.33				5.463	-	5.33			
		C-Zone	7.159	2.09	7.94	-			4.108	1.611	4.86			
12587-46-1	Gross Alpha [pci/L]	D-Zone	NA	NA	NA	3.2	0.09	0.56	NA	NA	NA	2.3	0.07	0.38
		Sunnyside	NA	NA	NA	1			NA	NA	NA			
		B-Zone	4.227	-	7.3				2.73	-	7.3			
12507 47 2	Corres Data Laci (1)	C-Zone	8.1	8.3	12	2.6	0.26	0.03	3.65	6.067	11	1.4	0.26	0.70
12587-47-2	Gross Beta [pci/L]	D-Zone	NA	NA	NA	2.6	0.36	0.83	NA	NA	NA	1.4	0.26	0.78
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
HAA5	HAA5, Total [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
ПААЗ	ΠΑΑ3, Total [μg/L]	D-Zone	ND	ND	ND	ND	IND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	750000	-	620000				690000	-	620000			
Hardt-1	Hardness, total as CaCO3 [µg/L]	C-Zone	700000	440000	550000	510000	290000	300000	540000	300000	540000	470000	280000	300000
		D-Zone	410000	410000	410000	-			410000	410000	410000			
		Sunnyside	270000	270000	270000				270000	270000	270000			
		B-Zone	ND	-	ND	-			ND	-	ND			
76-44-8	Heptachlor [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	,	D-Zone	ND	ND	ND	-			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND ND	- ND	ND	1			ND ND	- ND	ND			
1024-57-3	Heptachlor Epoxide [µg/L]	C-Zone D-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
		Sunnyside	ND ND	ND ND	ND ND	-			ND ND	ND ND	ND ND			
		B-Zone	ND ND	- ND	ND ND				ND ND	- ND	ND ND			
		C-Zone	ND ND	ND	ND	1			ND	- ND	ND ND			
111-71-7	Heptanal [μg/L]	D-Zone	NA NA	NA NA	NA NA	ND	ND	ND	NA NA	NA NA	NA NA	ND	ND	ND
		Sunnyside	NA NA	NA NA	NA	1			NA	NA	NA			







Case   Constituent [Units]   Concentration within Capture   Zone   SM-4   SM-8   SM-9   SM-4   SM-9   SM-9   SM-4   SM-9   SM-9   SM-4   SM-9   SM-9   SM-9   SM-4   SM-9   SM	SM-4  0.0002  ND  ND	Concentration Wells SM-8 0.003 ND	SM-9
118-74-1   Hexachlorobenzene [μg/L]   B-Zone   ND   -     0.014     0.009   ND     0.0002   0.0039   0.0003     0.0014     0.07   ND   ND   ND   ND   ND   ND   ND   N	0.0002 ND	0.003 ND	SM-9 0.0003 ND
Hexachlorobenzene [μg/L]	0.0002 ND	0.003 ND ND	0.0003 ND
Hexachlorobenzene [µg/L]	ND ND	ND ND	ND
Na   Na   Na   Na   Na   Na   Na   Na	ND ND	ND ND	ND
D-Zone   NA   NA   NA   NA   NA   NA   NA   N	ND ND	ND ND	ND
Hexachlorobutadiene [μg/L]   B-Zone   ND   -   ND   ND   ND   ND   ND   ND	ND	ND	
R7-68-3   Hexachlorobutadiene [μg/L]	ND	ND	
ND   ND   ND   ND   ND   ND   ND   ND	ND	ND	
D-Zone   ND   ND   ND   ND   ND   ND   ND   N	ND	ND	
Hexachlorocyclopentadiene [μg/L]   B-Zone   ND   -   ND   ND   ND   ND   ND   ND			ND
T7-47-4   Hexachlorocyclopentadiene [μg/L]   C-Zone   ND   ND   ND   ND   ND   ND   ND   N			ND
Hexachlorocyclopentadiene [μg/L]			ND
D-Zone   ND   ND   ND   ND   ND   ND   ND   N			
Hexachloroethane [μg/L]   B-Zone   ND   -   ND   ND   ND   ND   ND   ND	ND	ND	
Hexachloroethane [μg/L]	ND	ND	
D-Zone   ND   ND   ND   ND   ND   ND   ND   N	ND	ND	
D-Zone   ND   ND   ND   ND   ND   ND   ND   N		עוו ווע	ND
I R-7000 I ND I - I ND I I I I ND I I ND I			
66-25-1 Hexanal [μg/L] C-Zone ND	ND	ND	ND
D-Zone NA NA NA NA NA NA NA NA			
Sunnyside NA			
B-Zone ND - ND ND - ND ND - ND			
13252-13-6 HFPO-DA [μg/L] C-Zone ND	ND	ND	ND
D-Zone ND ND ND ND ND			
Sunnyside NA			
B-Zone ND - ND ND - ND			
2691-41-0 HMX [μg/L] C-Zone ND	ND	ND	ND
D-Zone NA			
B-Zone ND - ND			
302-01-2	ND	ND	ND
Sunnyside NA			
B-Zone 0.0048 - ND 0.0022 - ND			
C-7000 0.0023 0.0048 0.00066			
15687-27-1   Ibuprofen [μg/L]   C-201e 0.0023 0.0048 0.00000   0.0018   0.00021   4Ε-05   NA NA NA	0.0008	0.0001	3 4E-05
Sunnyside NA NA NA NA NA NA NA			
B-Zone ND - ND ND - ND			
C-Zone ND			
193-39-5 Indeno[1,2,3-cd]pyrene [μg/L]	ND	ND	ND
Sunnyside ND			
B-Zone 19 - 0.48 10 - 0.48			
C-7one 12 76 82 7 49 46			
20461-54-5   Iodide [μg/L]   C Zone   12	4.2	2.1	0.28
Sunnyside NA NA NA NA NA NA NA			







				Us	ing UCL95 fo	or Monitoring V	Well Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [[Inite]	Lover	Concent	ration within	Capture	Estimated C	oncentrations fo	or Production	Concent	ration within	Capture	Estimated Co	oncentrations fo	or Production
CASKIN	Constituent [Units]	Layer		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
74 00 4	la damathana [i.g./l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
74-88-4	lodomethane [μg/L]	D-Zone	ND	ND	ND	שא	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
73334-07-3	la promida (um/l.)	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
73334-07-3	lopromide [µg/L]	D-Zone	NA	NA	NA	שא	ND	ND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	10	-	ND				4.4	-	ND			
7439-89-6	lvon [uo/l]	C-Zone	3	2600	ND	7.2	120	8.7	3	950	ND	5.5	50	8.7
7439-09-0	lron [μg/L]	D-Zone	1.1	1.1	1.1	1.2	120	0.7	1.1	1.1	1.1	5.5	50	0.7
		Sunnyside	10	10	10				10	10	10			
		B-Zone	ND	-	ND				ND	-	ND			
78-59-1	Jaanharana [i.e./l.]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
70-39-1	Isophorone [μg/L]	D-Zone	ND	ND	ND	שא	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
67-63-0	Isanranyi Alsahal Iya/II	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
07-03-0	lsopropyl Alcohol [µg/L]	D-Zone	NA	NA	NA	ND	ND	IND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
98-82-8	Isanrandhanzana [ug/l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
90-02-0	Isopropylbenzene [µg/L]	D-Zone	ND	ND	ND	ND	ND	IND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.1	-	1.2				0.0843	-	1.2			
7439-92-1	Lead, Total [μg/L]	C-Zone	0.074	0.38	0.14	0.043	0.016	0.03	0.0727	0.165	0.14	0.038	0.0071	0.03
1439-92-1	Lead, Total [µg/L]	D-Zone	NA	NA	NA	0.043	0.016	0.03	NA	NA	NA	0.036	0.0071	0.03
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	_	ND			
330-55-2	Linuron [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
330 33 2	Linuion [μg/ L]	D-Zone	NA	NA	NA	I ND	ND	ND	NA	NA	NA	IND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
7439-93-2	Lithium [µg/L]	C-Zone	ND	25	ND	ND	1.1	ND	ND	18	ND	ND	0.76	ND
1433-33-2	Επιπατή [μg/ Ε]	D-Zone	NA	NA	NA	I ND	1.1	ND	NA	NA	NA	IND	0.70	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
179601-23-1	m,p-Xylene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
173001 23-1	m,p Ayiene [µg/L]	D-Zone	ND	ND	ND		IND	IND	ND	ND	ND	140	IND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	96000	-	62000				80000	-	62000			
7439-95-4	Magnesium [µg/L]	C-Zone	78000	51000	56000	58000	26000	27000	61000	36000	55000	50000	25000	27000
1733-33-4	iviagnesium [µg/L]	D-Zone	42000	42000	42000	30000	20000	27000	42000	42000	42000	30000	23000	27000
		Sunnyside	24000	24000	24000				24000	24000	24000			







				Usi	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ng Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	r Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	289.1	-	57				165	-	57			
7439-96-5	Manganoso Total Jug /L1	C-Zone	120	110	85	107.2	4.8	6	78.75	47.67	55.5	62.9	2.1	4.3
7459-90-5	Manganese, Total [μg/L]	D-Zone	NA	NA	NA	107.2	4.0	8	NA	NA	NA	02.9	2.1	4.5
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
2355-31-9	MeFOSAA [μg/L]	C-Zone	0.0013	ND	ND	0.0002	ND	ND	0.0013	ND	ND	0.0002	ND	ND
2333 31 3	ίνιει Ο ΣΑΑ [μg/.Ε]	D-Zone	ND	ND	ND	0.0002	ND	ND	ND	ND	ND	0.0002	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
57-53-4	Meprobamate [µg/L]	C-Zone	0.0022	ND	ND	0.0004	ND	ND	0.0014	ND	ND	0.0002	ND	ND
37 33 4	Meprobarilate [µg/1]	D-Zone	NA	NA	NA	0.0004	IND	ND	NA	NA	NA	0.0002	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.022	-	ND				0.022	-	ND	-		
7439-97-6	Mercury [µg/L]	C-Zone	ND	0.025	ND	0.017	0.019	0.018	ND	0.025	ND	0.017	0.019	0.018
7-335 37 0	(Well-edity [µg/-1]	D-Zone	0.023	0.023	0.023	0.017	0.013	0.010	0.021	0.021	0.021	0.017	0.013	0.010
		Sunnyside	0.019	0.019	0.019				0.019	0.019	0.019			
		B-Zone	ND	-	ND				ND	-	ND			
126-98-7	Methacrylonitrile [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
120 30 7	Methody Smalle [µg/ 2]	D-Zone	ND	ND	ND	110		112	ND	ND	ND		115	115
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.00024	-	ND				0.00015	-	ND	- -		
76-99-3	Methadone [µg/L]	C-Zone	0.00022	0.000063	ND	0.0001	2.7E-06	ND	0.00013	0.000063	ND	7E-05	2.7E-06	ND
		D-Zone	NA	NA	NA				NA	NA	NA	-		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND	-		
67-56-1	Methanol [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	11 37 3	D-Zone	NA	NA	NA				NA	NA	NA	-		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
2032-65-7	Methiocarb [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	- ND	ND	-		
16752-77-5	Methomyl [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	· -	D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	- NID	ND	-		
72-43-5	Methoxychlor [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	- ND	ND	-		
78-98-8	Methyl Glyoxal [µg/L]	C-Zone	ND	ND NA	ND NA	ND	ND	ND	ND NA	ND NA	ND NA	ND	ND	ND
		D-Zone	NA	NA NA	NA NA				NA NA	NA	NA NA	-		
		Sunnyside	NA	NA	NA		]		NA	NA	NA			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	r Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	0.13	-	0.083				0.13	-	0.083			
00.63.6	A	C-Zone	ND	ND	ND	0.020	NB	0.0045	ND	ND	ND	0.000	ND	0.0015
80-62-6	Methyl Methacrylate [µg/L]	D-Zone	ND	ND	ND	0.039	ND	0.0015	ND	ND	ND	0.039	ND	0.0015
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.53	-	ND				0.53	-	ND			
1634-04-4	Mothyl tort butyl other (MTPE) [ug/l]	C-Zone	ND	3.8	ND	0.16	0.17	ND	ND	1.8	ND	0.16	0.077	ND
1634-04-4	Methyl tert-butyl ether (MTBE) [μg/L]	D-Zone	ND	ND	ND	0.16	0.17	ND	ND	ND	ND	0.16	0.077	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	34	-	ND				26	-	ND			
MBAS	Mathylana blua activa substances (MDAC) [ug/l]	C-Zone	19	38	20	52	79	76	19	27	20	47	78	75
IVIDAS	Methylene blue active substances (MBAS) [µg/L]	D-Zone	52	52	52	52	79	76	42	42	42	47	70	/5
		Sunnyside	83	83	83				83	83	83			
		B-Zone	ND	-	ND				ND	-	ND			
75-09-2	Mothydono Chlorido Iva (L1	C-Zone	ND	0.81	ND	0.4	0.72	0.66	ND	0.64	ND	0.37	0.71	0.65
75-09-2	Methylene Chloride [μg/L]	D-Zone	0.85	0.85	0.85	0.4	0.72	0.00	0.68	0.68	0.68	0.57	0.71	0.65
		Sunnyside	0.71	0.71	0.71				0.71	0.71	0.71			
		B-Zone	ND	-	ND				ND	-	ND			
51218-45-2	Motologblev [ug/l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
31210-43-2	Metolachlor [µg/L]	D-Zone	ND	ND	ND	ND	IND	IND	ND	ND	ND	] ND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	_	ND				ND	-	ND			
171118-09-5	Metolachlor ESA [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
171110-09-5	Wetolacillor L3A [μg/L]	D-Zone	NA	NA	NA	ND	IND	IND	NA	NA	NA	IND	ND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
152019-73-3	Metolachlor OA [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
132013-73-3	Wetolachiol OA [μg/ L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA		ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
21087-64-9	Metribuzin [μg/L]	C-Zone	ND	0.079	ND	ND	0.0034	ND	ND	0.079	ND	ND	0.0034	ND
21007 04 3	(μg/ ε]	D-Zone	ND	ND	ND	110	0.0054	110	ND	ND	ND		0.0034	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
2385-85-5	Mirex [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
2303 03 3	ινιι ελ [μg/ -]	D-Zone	ND	ND	ND	115		115	ND	ND	ND		115	
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND	1		
2212-67-1	Molinate [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND			. 15	ND	ND	ND	1		1,5
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	14	-	12				11	-	12	1		
7439-98-7	Molybdenum [μg/L]	C-Zone	60	46	16	14	2	1.2	27	32	12	7.7	1.4	0.94
55 50 .		D-Zone	NA	NA	NA				NA	NA	NA			3.3 .
		Sunnyside	NA	NA	NA				NA	NA	NA			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	· Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concentr	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production	Concentr	ation within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
70.00.3		C-Zone	ND	ND	ND	ND	NB	ND	ND	ND	ND	ND	ND	ND
79-08-3	Monobromoacetic acid [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	7.7	1	100				6	-	100			
10599-90-3	Manachlaramina [ug/l]	C-Zone	9	8	10	3.8	0.35	2.4	4.7	6.5	7	2.6	0.28	2.2
10599-90-3	Monochloramine [µg/L]	D-Zone	NA	NA	NA	3.0	0.35	2.4	NA	NA	NA	2.0	0.26	2.2
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
60-34-4	Monomethylhydrazine [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
00-34-4	Monomethymydrazme [μg/L]	D-Zone	NA	NA	NA	ND	IND	IND	NA	NA	NA	ND	ND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.0011	-	0.00096				0.00084	-	0.00096			
57-27-2	Morphine [µg/L]	C-Zone	0.0018	0.0018	0.00094	0.0006	7.8E-05	7E-05	0.0017	0.0016	0.00094	0.0005	6.7E-05	7E-05
31-21-2	Могрине [µg/L]	D-Zone	NA	NA	NA	0.0000	7.0L-03	712-03	NA	NA	NA	0.0003	0.7L-03	76-05
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
91-20-3	Naphthalene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
31 20 3	ιναρπιαίστε [μg/ 2]	D-Zone	ND	ND	ND	ND	110	110	ND	ND	ND	110	ND	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.0029	-	ND				0.00095	-	ND			
22204-53-1	Naproxen [µg/L]	C-Zone	ND	ND	ND	0.0009	ND	ND	ND	ND	ND	0.0003	ND	ND
		D-Zone	NA	NA	NA				NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
104-51-8	n-Butylbenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 -	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			+
		B-Zone	200	-	7.2				71.15	-	7.2			
7440-02-0	Nickel, Total [μg/L]	C-Zone	84	3.6	190	74.3	0.16	11.3	45.28	2.05	138	29	0.089	8.2
		D-Zone	NA	NA	NA				NA	NA NA	NA			
		Sunnyside	NA 12.74	NA	NA 10				NA C 05	NA	NA 10			
		B-Zone	12.74 9.7	- 4.2	10				6.85	1 100	10			
14797-55-8	Nitrate as N [mg/L]	C-Zone D-Zone	0.83	4.2 0.83	6.4 0.83	5.6	0.23	0.6	6.033 0.83	1.185 0.83	4.5 0.83	3.2	0.1	0.49
		Sunnyside	0.63 ND	0.63 ND	ND				ND	ND	ND			
		B-Zone	99.23	- -	ND ND				65.83	-	ND			
		C-Zone	76	- ND	58				74	ND	58			
14797-65-0	Nitrite as N [µg/L]	D-Zone	50	50	50	51.7	2.8	6.2	50	50	50	41.3	2.8	6.2
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND ND				ND	-	ND ND			
		C-Zone	ND	ND	ND				ND	ND	ND			
98-95-3	Nitrobenzene [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND ND				ND	ND ND	ND			
		Surinyside	ND	ND	טאו			1	ND	טאו	שוו			







				Us	ing UCL95 f	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	/ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	r Production	Concent	ration withir Zone	Capture	Estimated C	oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND	1			ND	ND	ND			
55-18-5	N-Nitrosodiethylamine (NDEA) [ng/L]	D-Zone	1.1	1.1	1.1	0.2	0.061	0.062	1.1	1.1	1.1	0.2	0.061	0.062
		Sunnyside	ND	ND	ND	1			ND	ND	ND	1		
		B-Zone	ND	-	ND				ND	-	ND			
62.75.0	NAME OF THE PARTY	C-Zone	0.0018	ND	ND	0.000	ND	NB	0.0018	ND	ND	0.000	ND	ND
62-75-9	N-Nitrosodimethylamine [µg/L]	D-Zone	ND	ND	ND	0.0003	ND	ND	ND	ND	ND	0.0003	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
024.46.2	Alter III and the Control	C-Zone	ND	ND	ND	1	ND	NB	ND	ND	ND		ND	ND
924-16-3	n-Nitrosodi-n-butylamine [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
C21 C4 7	n Nitraca di manana la mina five // 1	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
621-64-7	n-Nitrosodi-n-propylamine [µg/L]	D-Zone	ND	ND	ND	שוא	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
86-30-6	n Nitro co dinhamulamina [um/l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
86-30-6	n-Nitrosodiphenylamine [μg/L]	D-Zone	ND	ND	ND	שוא	ND	שא	ND	ND	ND	שא	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
10595-95-6	N-Nitrosomethylethylamine [µg/L]	C-Zone	ND	ND	ND	ND ND	ND	ND	ND	ND	ND	ND	ND	ND
10393-93-0	iv-ivitiosometriyletriylarilile [µg/L]	D-Zone	ND	ND	ND		IND	ND	ND	ND	ND	IND	IND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.031	-	0.0017				0.016	-	0.0017			
59-89-2	n-Nitrosomorpholine [µg/L]	C-Zone	0.013	0.0032	0.0015	0.012	0.00014	0.0001	0.0078	0.0014	0.0012	0.0062	6.2E-05	0.0001
33 03 2	π wittosomorpholine [μg/ L]	D-Zone	ND	ND	ND	0.012	0.00014	0.0001	ND	ND	ND	0.0002	0.21-03	0.0001
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
100-75-4	n-Nitrosopiperidine [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
100 73 4	π τια σσοριμοπαιπο [μg/ 2]	D-Zone	ND	ND	ND		110	110	ND	ND	ND		110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND	_		
930-55-2	n-Nitrosopyrrolidine [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND	1			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	16	-	10	_			6.74	-	10	-		
NN	NO2+NO3 as N [mg/L]	C-Zone	9.536	2.526	6.4	6.9	1	1.4	6.025	1.225	4.5	3.6	0.95	1.3
		D-Zone	1.15	1.15	1.15	_			1.15	1.15	1.15			
		Sunnyside	0.92	0.92	0.92				0.92	0.92	0.92			
		B-Zone	0.41	-	ND	1			0.095	-	ND	-		
25154-52-3	Nonylphenol [μg/L]	C-Zone	ND	4.4	ND	0.12	0.19	ND	ND	3.5	ND	0.029	0.15	ND
	7 1 60 3	D-Zone	NA	NA	NA	1			NA	NA	NA	-		
		Sunnyside	NA	NA	NA			]	NA	NA	NA	1		







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ing Mean for	· Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concentr	ation within Zone	Capture	Estimated C	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	0.048	-	ND				0.048	-	ND			
		C-Zone	ND	3.2	ND				ND	2.3	ND			
20427-84-3	Nonylphenol diethoxylate [µg/L]	D-Zone	NA	NA	NA	0.014	0.14	ND	NA	NA	NA	0.014	0.1	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.049	-	ND				0.026	-	ND			
27006 26 2		C-Zone	0.045	2.3	0.013		0.4	0.000	0.032	1.8	0.013	0.010	0.070	0.000
27986-36-3	Nonylphenol monoethoxylate [µg/L]	D-Zone	NA	NA	NA	0.022	0.1	0.0008	NA	NA	NA	0.013	0.078	0.0008
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
102 (5.1	n Dramillandana (i.e./l.)	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
103-65-1	n-Propylbenzene [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	=	ND			
136777-61-2	o&p-Xylene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
130///-01-2	οαρ-λуιεπε [μg/L]	D-Zone	ND	ND	ND	IND	IND	ND	ND	ND	ND	ND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	NA	-	NA				NA	-	NA			
111-66-0	Octene-1 [µg/L]	C-Zone	25	NA	NA	4.2	NA	NA	25	NA	NA	4.2	NA	NA
111 00 0	σετεπε τ [μg/ ε]	D-Zone	NA	NA	NA	7.2	IN/A	IVA	NA	NA	NA	7.2	l IVA	IVA
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	1500	-	ND	_			1400	-	ND			
OILGREASE	Oil & Grease (HEM) [µg/L]	C-Zone	ND	ND	ND	450	ND	ND	ND	ND	ND	420	ND	ND
	o o. o. o o ( , t[mg, -]	D-Zone	NA	NA	NA	-			NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	78	-	150	-			61	-	150			
14265-44-2	o-Phosphate as P [μg/L]	C-Zone	83	95	68	37	4.1	6.7	63	82	65	29	3.5	6.5
	,	D-Zone	NA	NA	NA	-			NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND	-			ND	- ND	ND			
23135-22-0	Oxamyl [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND ND	ND ND	ND ND	-			ND ND	ND ND	ND ND			
		Sunnyside B-Zone	0.012	-	0.006				0.0065	- ND	0.006			
		C-Zone	0.012	0.029	0.006	1			0.0063	0.02	0.008			
131-57-7	Oxybenzone [µg/L]	D-Zone	0.0032 NA	0.029 NA	0.0094 NA	0.0042	0.0013	0.0007	0.0017 NA	NA	0.0088 NA	0.0023	0.00087	0.0006
		Sunnyside	NA NA	NA NA	NA NA	-			NA NA	NA NA	NA NA			
		B-Zone	ND ND	INA	ND				ND ND	- INA	ND ND			
		C-Zone	ND ND	0.31	ND ND	1			ND ND	0.31	ND ND			
95-47-6	o-Xylene [μg/L]	D-Zone	ND ND	ND	ND ND	ND	0.013	ND	ND ND	ND	ND ND	ND	0.013	ND
		Sunnyside	ND ND	ND	ND	-			ND ND	ND ND	ND			
		B-Zone	ND ND	-	ND				ND	-	ND			
		C-Zone	ND ND	ND	ND	1			ND	ND	ND			
4685-14-7	Paraquat [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND	1			ND	ND	ND			







				Us	sing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	or Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
100004 64 0	DDDE 100 for #11	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	l ND	ND	ND
189084-64-8	PBDE-100 [µg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
182677-30-1	PBDE-138 [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
102077-30-1	ΡΒ <b>ΔΕ-</b> 130 [μg/L]	D-Zone	NA	NA	NA	IND	ND	IND	NA	NA	NA	IND	IND	IND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
68631-49-2	PBDE-153 [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
00031-49-2	νωυ <u>ς</u> -155 [μg/ε]	D-Zone	NA	NA	NA	IND	ND	IND	NA	NA	NA		IND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
207122-15-4	PBDE-154 [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
20/122 13 4	1 υυς 134 [μg/ε]	D-Zone	NA	NA	NA	I ND	ND	ND	NA	NA	NA		ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
147217-75-2	PBDE-17 [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
147217 732	1 2 2 17 [μg/ ε]	D-Zone	NA	NA	NA		145	145	NA	NA	NA	1	110	110
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
41318-75-6	PBDE-28 [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	: 22 = 20 [#g/ =]	D-Zone	NA	NA	NA				NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND	_		
5436-43-1	PBDE-47 [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	LF 3, 1	D-Zone	NA	NA	NA				NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
243982-82-3	PBDE-49 [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	-, 5	D-Zone	NA NA	NA	NA	_			NA	NA	NA	_		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	- ND	ND	_			ND	- ND	ND	_		
182346-21-0	PBDE-85 [μg/L]	C-Zone	ND NA	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	NA NA	NA	NA	-			NA	NA	NA	_		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	- ND	ND	_			ND	- ND	ND	-		
60348-60-9	PBDE-99 [µg/L]	C-Zone	ND	ND NA	ND NA	ND	ND	ND	ND NA	ND NA	ND NA	ND	ND	ND
		D-Zone	NA	NA NA	NA NA	1			NA NA	NA NA	NA	+		
		Sunnyside	NA ND	NA	NA					NA	NA			+
		B-Zone	ND ND	- ND	ND	1			ND ND	- ND	ND	1		
76-01-7	Pentachloroethane [µg/L]	C-Zone	ND ND	ND ND	ND ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
		D-Zone			_	1								
		Sunnyside	ND	ND	ND				ND	ND	ND			







				Usi	ing UCL95 fo	r Monitoring V	Vell Data*			Usi	ing Mean fo	r Monitoring W	ell Data**	
			Concent	ration within			oncentrations fo	or Production	Concent	ration within			oncentrations fo	r Production
CASRN	Constituent [Units]	Layer		Zone	omp om o		Wells			Zone	p		Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
82-68-8	Pentachloronitrobenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
02 00 0	r entachioronitrobenzene [μg/ ε]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
87-86-5	Pentachlorophenol [µg/L]	C-Zone	ND	0.43	ND	ND	0.019	ND	ND	0.29	ND	ND	0.013	ND
07 00 3		D-Zone	ND	ND	ND	ND	0.013	110	ND	ND	ND	110	0.013	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
110-62-3	Pentanal [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	. 5a. [h-3, -]	D-Zone	NA	NA	NA				NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	6.7	-	3.3				1.817	-	3.3			
14797-73-0	Perchlorate [μg/L]	C-Zone	3.3	2	2	2.6	0.087	0.18	2.367	2	1.8	0.94	0.087	0.16
	11 37 1	D-Zone	NA	NA	NA				NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.0056	-	0.0036				0.0041	-	0.0036			
375-73-5	PFBS [μg/L]	C-Zone	0.0031	0.0014	0.0018	0.0022	6.1E-05	0.0002	0.0021	0.0014	0.0018	0.0016	6.1E-05	0.0002
	5 -	D-Zone	ND	ND	ND NA				ND	ND	ND			
		Sunnyside	NA 0.0020	NA	NA				NA 0.0015	NA	NA			
		B-Zone	0.0028	- ND	ND 0.0031				0.0015	- ND	ND 0.0010			
335-76-2	PFDA [µg/L]	C-Zone	0.0062 ND	ND ND	0.0021 ND	0.0019	ND	0.0001	0.0045 ND	ND ND	0.0019 ND	0.0012	ND	0.0001
		D-Zone	NA NA		NA NA				NA NA	NA NA	NA NA			
		Sunnyside B-Zone	0.00073	NA -	NA ND				0.00061	INA -	ND ND			
		C-Zone	0.00073	0.0018	0.00056				0.00049	0.0017	0.00056			
307-55-1	PFDoA [μg/L]	D-Zone	ND	0.0018 ND	ND	0.0003	7.8E-05	3E-05	0.00049 ND	0.0017 ND	ND	0.0003	7.5E-05	3E-05
		Sunnyside	NA NA	NA NA	NA NA				NA NA	NA NA	NA NA			
		B-Zone	0.0032	-	0.0016				0.0022	-	0.0016			
		C-Zone	0.0032	0.002	0.00098				0.0022	0.0014	0.00016			
375-85-9	PFHpA [μg/L]	D-Zone	0.0015	0.0055	0.0055	0.0022	0.00039	0.0004	0.0012	0.0037	0.0003	0.0015	0.00027	0.0003
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.012	-	0.0053				0.0078	-	0.0053			
		C-Zone	0.0054	0.0025	0.0021				0.0031	0.0016	0.0015			
307-24-4	PFHxA [μg/L]	D-Zone	0.0007	0.0007	0.0007	0.0047	0.00015	0.0003	0.0007	0.0007	0.0007	0.003	0.00011	0.0002
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.0063	-	0.0035				0.0042	-	0.0035			
		C-Zone	0.0032	0.0013	0.0025				0.0024	0.0013	0.0021			
355-46-4	PFHxS [μg/L]	D-Zone	ND	ND	ND	0.0024	5.6E-05	0.0002	ND	ND	ND	0.0017	5.4E-05	0.0002
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.0016	-	ND				0.0014	-	ND			
275 05 4	DENIA C. (1)	C-Zone	0.0047	ND	ND	0.0013		N.S	0.0026	ND	ND	0.0000	NE	
375-95-1	PFNA [μg/L]	D-Zone	ND	ND	ND	0.0013	ND	ND	ND	ND	ND	0.0009	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ng Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	or Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	3.135	-	2				1.902	-	2			
225 67 4	DEO.A. (	C-Zone	0.971	0.89	1.4	4.4	0.020	0.12	0.668	0.655	1.4	0.60	0.000	0.10
335-67-1	PFOA [ng/L]	D-Zone	ND	ND	ND	1.1	0.039	0.12	ND	ND	ND	0.68	0.028	0.12
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	4.2				ND	-	4.2			
1763-23-1	DEOC [na/L]	C-Zone	ND	ND	ND	ND	ND	0.075	ND	ND	ND	ND	ND	0.075
1703-23-1	PFOS [ng/L]	D-Zone	ND	ND	ND	ND	ND	0.075	ND	ND	ND	IND	ND	0.075
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
376-06-7	PFTeDA [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
370-00-7	ΓΙ ΤΕ <b>Ο</b> Α [μg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	I ND	IND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
72629-94-8	PFTrDA [μg/L]	C-Zone	ND	0.00041	ND	ND	1.8E-05	ND	ND	0.00041	ND	ND	1.8E-05	ND
72023 34 0	11110Α [μg/ ε]	D-Zone	ND	ND	ND	ND	1.02 03	IND	ND	ND	ND	I ND	1.02 03	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.0028	-	ND				0.0022	-	ND			
2058-94-8	PFUnA [μg/L]	C-Zone	0.0024	0.00078	0.0031	0.0013	3.4E-05	0.0002	0.002	0.00078	0.0024	0.001	3.4E-05	0.0001
2030 34 0	11 Οτην [μβ/ 2]	D-Zone	ND	ND	ND	0.0013	3.42 03	0.0002	ND	ND	ND	0.001	3.42 03	0.0001
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	NA	-	NA				NA	-	NA			
TPHd	PHC AS DIESEL FUEL [µg/L]	C-Zone	NA	100	NA	NA	4.3	NA	NA	67	NA	NA NA	2.9	NA
	: e , . e	D-Zone	NA	NA	NA		5		NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND	-		
85-01-8	Phenanthrene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
108-95-2	Phenol [μg/L]	C-Zone	ND	1.2	ND	ND	0.052	ND	ND	1	ND	ND	0.045	ND
	5· ·	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- ND	ND				ND	-	ND	-		
57-41-0	Phenytoin (Dilantin) [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND NA	ND	ND	ND	ND	ND
		D-Zone	NA	NA	NA				NA NA	NA	NA	-		
		Sunnyside	NA 250	NA	NA				NA 200	NA	NA 600			
		B-Zone	250	200	680				200	100	680	-		
7723-14-0	Phosphorus [µg/L]	C-Zone	720 NA	300	730	200	13	55	380	180	630	120	7.7	49
		D-Zone	NA NA	NA NA	NA NA				NA NA	NA NA	NA NA	-		
		Sunnyside	NA o a	NA	NA NA				NA 4.8	NA	NA NA			
		B-Zone	8.3 4.6	- 15	NA NA				4.8 4.6	8.6	NA NA	-		
85-44-9	Phthalic Anhydride [µg/L]	C-Zone			NA NA	3.3	0.65	NA			NA NA	2.2	0.37	NA
		D-Zone	NA NA	NA NA	NA NA				NA NA	NA NA	NA NA	-		
		Sunnyside	NA	NA	NA				NA	NA	NA			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
			Concent	ration within	Capture	Estimated C	oncentrations fo	or Production	Concent	ration within			oncentrations fo	or Production
CASRN	Constituent [Units]	Layer		Zone			Wells			Zone	p		Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	=	ND				ND	-	ND			
1918-02-1	Picloram [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1910-02-1	Ficiorani [μg/L]	D-Zone	ND	ND	ND	I ND	IND	IND	ND	ND	ND	IND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
99-87-6	p-lsopropyltoluene [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
33 01 0	ρ ισοριοργιτοιατίτε [μg/ ε]	D-Zone	ND	ND	ND	l ND	ND	I II	ND	ND	ND	110	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	4200	-	3300				3500	-	3300			
7440-09-7	Potassium [μg/L]	C-Zone	5600	10000	3100	4600	5200	4900	4300	5900	2900	4200	5000	4800
7440 03 7	Ι οτασσιατίτ [μg/ ε]	D-Zone	3600	3600	3600	-1000	3200	4500	3500	3500	3500	4200	3000	4000
		Sunnyside	5100	5100	5100				5100	5100	5100			
		B-Zone	ND	-	ND				ND	-	ND			
55268-74-1	Praziquantel [µg/L]	C-Zone	ND	0.0066	ND	ND	0.00029	ND	ND	0.0038	ND	ND	0.00016	ND
33233 7		D-Zone	NA	NA	NA		0.00025		NA	NA	NA		0.000.0	
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.0029	-	0.0018				0.0018	-	0.0018			
125-33-7	Primidone [µg/L]	C-Zone	ND	ND	ND	0.0009	ND	3E-05	ND	ND	ND	0.0005	ND	3E-05
		D-Zone	NA	NA	NA			02.00	NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
57-83-0	Progesterone [μg/L]	C-Zone	ND	0.00033	ND	ND	1.4E-05	ND	ND	0.00033	ND	ND	1.4E-05	ND
	3 1 3 3	D-Zone	NA	NA	NA				NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			+
		B-Zone	ND	-	ND				ND	-	ND			
1610-18-0	Prometon [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	., 3, -	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
7287-19-6	Prometryn [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND	-			ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			
	December (2 Chloro et (1 Marthelath ) et about la	B-Zone	ND	- ND	ND	-			ND ND	- ND	ND	-		
1918-16-7	Propachlor (2-Chloro-n-(1-Methylethy)-n-phenylac	C-Zone	ND	ND	ND	ND	ND	ND	ND ND	ND	ND	ND	ND	ND
	[µg/L]	D-Zone	ND ND	ND ND	ND ND	-			ND ND	ND ND	ND ND	-		
		Sunnyside B-Zone	1.2	-	ND						ND			
		C-Zone	1.4	- ND	ND ND	-			1.2	- ND	ND ND			
123-38-6	Propanal [μg/L]	D-Zone	NA	NA NA	NA NA	0.6	ND	ND	NA	NA NA	NA NA	0.6	ND	ND
		Sunnyside	NA NA	NA NA	NA NA	1			NA NA	NA NA	NA NA			
		B-Zone	ND ND	INA -	ND				ND ND	INA -	ND			
		C-Zone	ND ND	- ND	ND	1			ND	- ND	ND			
114-26-1	Propoxur (Baygon) [μg/L]	D-Zone	ND ND	ND	ND	ND	ND	ND	ND	ND ND	ND	ND	ND	ND
		Sunnyside	ND ND	ND ND	ND	1			ND	ND ND	ND			
		Surinysiae	ND	ND	ND				ואט	ND	ND			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Usi	ng Mean foi	r Monitoring W	/ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within	Capture	Estimated C	oncentrations fo	r Production	Concent	ation within	Capture	Estimated C	oncentrations fo	r Production
CASKIN	Constituent [onits]	Layei		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
129-00-0	Pyrene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
123 00 0	r yrene [μg/ L]	D-Zone	ND	ND	ND		110	IND	ND	ND	ND	110	110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	  -			ND	-	ND			
110-86-1	Pyridine [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	y tr 5/ 1	D-Zone	ND	ND	ND	-			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.003	- 0.000	ND 0.0015	-			0.002	- 0.006	ND 0.0013			
91-22-5	Quinoline [µg/L]	C-Zone	0.0019	0.006	0.0015	0.0012	0.00026	9E-05	0.0017	0.006	0.0013	0.0009	0.00026	7E-05
		D-Zone	NA	NA NA	NA NA	-			NA NA	NA NA	NA NA			
		Sunnyside B-Zone	NA 0.205	NA -	0.12				NA 0.041	NA -	0.12			
		C-Zone	0.203	0.461	0.12	-			0.041	0.204	0.12			
13982-63-3	Radium 226 [pci/L]	D-Zone	0.323 NA	NA	NA	0.15	0.02	0.014	NA	NA	NA	0.048	0.0088	0.01
		Sunnyside	NA NA	NA NA	NA NA	-			NA	NA NA	NA			
		B-Zone	0.563	-	0.0865				0.317	-	0.0865			
		C-Zone	0.98	0.998	0.325	1			0.457	0.433	0.218			
15262-20-1	Radium 228 [pci/L]	D-Zone	NA	NA	NA	0.33	0.043	0.021	NA	NA	NA	0.17	0.019	0.014
		Sunnyside	NA	NA	NA	<u> </u>			NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
104.00		C-Zone	ND	ND	ND	1			ND	ND	ND			
121-82-4	RDX [μg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.58	-	0.45				0.39	-	0.45			
69-72-7	Caliculia Acid [ug/l]	C-Zone	0.59	0.74	0.36	0.27	0.032	0.029	0.29	0.41	0.34	0.17	0.018	0.028
09-72-7	Salicylic Acid [µg/L]	D-Zone	NA	NA	NA	0.27	0.032	0.029	NA	NA	NA	0.17	0.016	0.026
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
135-98-8	sec-Butylbenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
133 30 0	see batylisenzene [µg/ L]	D-Zone	ND	ND	ND		110	145	ND	ND	ND	, , ,	110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	5.247	-	3.2	-			3.795	-	3.2			
7782-49-2	Selenium, Total [μg/L]	C-Zone	3.4	0.72	3.8	2.1	0.031	0.28	2.725	0.347	2.85	1.6	0.015	0.22
		D-Zone	NA	NA	NA	-			NA	NA	NA			
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	34000	-	29000	-			32000	16000	29000			
7631-86-9	Silica [µg/L]	C-Zone D-Zone	35000	32000	35000	16000	1400	2600	34000	16000	34000	15000	700	2500
			NA NA	NA NA	NA NA	1			NA NA	NA NA	NA NA			
		Sunnyside B-Zone	ND ND	NA -	NA ND				NA ND	NA -	NA ND			
		C-Zone	ND ND	- ND	ND	1			ND ND	ND	ND ND			
7440-22-4	Silver [μg/L]	D-Zone	ND ND	ND ND	ND	ND	ND	ND	ND	ND ND	ND ND	ND	ND	ND
		Sunnyside	ND ND	ND ND	ND	-			ND	ND ND	ND ND			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	· Monitoring W	ell Data**	
CACRNI	Constitution III deal		Concent	ration within	Capture	Estimated Co	oncentrations fo	or Production	Concenti	ation within	Capture	Estimated Co	oncentrations fo	or Production
CASRN	Constituent [Units]	Layer		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
122-34-9	Simazina [ug/l]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
122-34-9	Simazine [µg/L]	D-Zone	ND	ND	ND	ND	IND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	100000	-	110000				84000	_	110000			
7440-23-5	Sodium [µg/L]	C-Zone	110000	120000	90000	78000	52000	53000	88000	89000	77000	68000	51000	52000
7440-23-3	30didiii [µg/L]	D-Zone	58000	58000	58000	70000	32000	33000	58000	58000	58000	00000	31000	32000
		Sunnyside	49000	49000	49000				49000	49000	49000			
		B-Zone	0.231	-	ND				0.0602	_	ND			
10098-97-2	Strontium-90 [pci/L]	C-Zone	0.461	0.428	0.045	0.25	0.27	0.25	0.207	0.143	0.0225	0.15	0.26	0.25
10090-91-2	Strontium-30 [pci/L]	D-Zone	NA	NA	NA	0.23	0.27	0.23	NA	NA	NA	0.13	0.20	0.23
		Sunnyside	0.282	0.282	0.282				0.282	0.282	0.282			
		B-Zone	ND	-	ND				ND	-	ND			
100-42-5	Styrene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
100-42-3	Styrene [µg/L]	D-Zone	ND	ND	ND	ND	IND	IND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	_	ND			
56038-13-2	Sucralose [µg/L]	C-Zone	0.0086	ND	ND	0.0014	ND	ND	0.0062	ND	ND	0.001	ND	ND
30030-13-2	Sucraiose [µg/L]	D-Zone	NA	NA	NA	0.0014	IND	IND	NA	NA	NA	0.001	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.00027	-	ND				0.00027	-	ND			
723-46-6	Sulfamethoxazole [µg/L]	C-Zone	ND	ND	ND	8E-05	ND	ND	ND	ND	ND	8E-05	ND	ND
725-40-0	Sunamethoxazole [µg/L]	D-Zone	NA	NA	NA	0L-03	IND	IND	NA	NA	NA	0L-03	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	300000	-	220000				270000	-	220000			
14808-79-8	Sulfate as SO4 [µg/L]	C-Zone	1600000	170000	220000	420000	100000	110000	680000	140000	210000	260000	88000	96000
14000 73 0	Surface as SO4 [µg/L]	D-Zone	180000	180000	180000	420000	100000	110000	180000	180000	180000	200000	00000	30000
		Sunnyside	96000	96000	96000				80000	80000	80000			
		B-Zone	ND	-	ND				ND	-	ND			
18496-25-8	Sulfide [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
10-130 23 0	Samae [pg/L]	D-Zone	NA	NA	NA	110	110	110	NA	NA	NA	110	ND	145
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.00055	-	ND				0.00055	_	ND			
115-96-8	TCEP [µg/L]	C-Zone	0.002	0.0006	ND	0.0005	2.6E-05	ND	0.0015	0.0006	ND	0.0004	2.6E-05	ND
113 30 0	[μg/ε]	D-Zone	NA	NA	NA	0.0003	2.02 03	110	NA	NA	NA	0.0004	2.02 03	145
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.0043	-	0.0013				0.0028	-	0.0013			
13674-84-5	TCPP [µg/L]	C-Zone	0.008	0.0069	0.0047	0.0026	0.0003	0.0003	0.003	0.0053	0.0028	0.0014	0.00023	0.0002
15074 04 5	τειτ [μg/ -]	D-Zone	NA	NA	NA	0.0020	0.0003	0.0003	NA	NA	NA	0.0014	0.00023	0.0002
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	0.0034	-	0.0026				0.0021	-	0.0026			
13674-87-8	TDCPP [µg/L]	C-Zone	0.0039	ND	0.0016	0.0017	ND	0.0001	0.0025	ND	0.0014	0.001	ND	0.0001
1307 4 07 0	12011 [μg/L]	D-Zone	NA	NA	NA	0.0017	IND	3.0001	NA	NA	NA	0.001	ND	0.0001
		Sunnyside	NA	NA	NA				NA	NA	NA			







				Us	ing UCL95 f	or Monitoring \	Well Data*			Usi	ng Mean fo	r Monitoring W	'ell Data**	
CACDN	Countition and Illustral		Concent	ration within	Capture	Estimated C	oncentrations fo	or Production	Concent	ration within	Capture	Estimated C	oncentrations fo	r Production
CASRN	Constituent [Units]	Layer		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	2.4	-	NA				1.7	-	NA			
TIC	Tentatively Identified Compounds [µg/L]	C-Zone	6.7	12	NA	1.8	0.51	NA	4.3	8.4	NA	1.2	0.36	NA
l lic	rentatively identified Compounds [µg/L]	D-Zone	NA	NA	NA	1.0	0.51	INA	NA	NA	NA	1.2	0.50	INA
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
5902-51-2	Terbacil [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
3302 31 2	rerbuen [µg/L]	D-Zone	ND	ND	ND		IND	IND	ND	ND	ND	, ND	IND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
13071-79-9	Terbufos [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
13071 73 3	τεισαίος [μg/ ε]	D-Zone	NA	NA	NA		140	145	NA	NA	NA		110	110
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
98-06-6	tert-Butylbenzene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND ND	ND	ND	ND	ND
	tert 2 aty. 2 en 2 en 2 [µg/ 2]	D-Zone	ND	ND	ND	1			ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	_			ND	-	ND			
994-05-8	Tertiary-amyl methyl ether [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
		D-Zone	ND	ND	ND	1			ND	ND	ND	-		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND	1			ND	-	ND	-		
75-65-0	Tertiary-butyl alcohol [µg/L]	C-Zone	ND	8.1	ND	ND	0.35	ND	ND	2	ND	ND	0.088	ND
	, ,	D-Zone	ND	ND	ND	-			ND	ND	ND	-		
		Sunnyside	ND 0.00025	ND	ND				ND 0.00033	ND	ND			
		B-Zone C-Zone	0.00025 ND	0.00034	ND ND	+			0.00022 ND	0.00034	ND ND	-		
58-22-0	Testosterone [µg/L]		NA NA	0.00034 NA	NA NA	8E-05	1.5E-05	ND	NA NA	0.00034 NA	NA NA	7E-05	1.5E-05	ND
		D-Zone	NA NA	NA NA	NA NA	+			NA NA	NA NA	NA NA	-		
		Sunnyside B-Zone	21.57	INA -	48.52				16.73	- INA	40.28			
		C-Zone	126.8	17.82	11.48	-			65.31	7.861	7.782	-		
127-18-4	Tetrachloroethene ( <mark>PCE)</mark> [µg/L]	D-Zone	ND	ND	ND	27.7	0.77	1.5	ND	ND	ND	16	0.34	1.2
		Sunnyside	ND ND	ND	ND ND				ND	ND ND	ND	-		
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND	1			ND	ND	ND	-		
109-99-9	Tetrahydrofuran [µg/L]	D-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND	1			ND	ND	ND	1		
479-45-8	Tetryl [µg/L]	D-Zone	NA	NA	NA	ND	ND	ND	NA	NA	NA	ND	ND	ND
		Sunnyside	NA	NA	NA	1			NA	NA	NA	1		
		B-Zone	0.019	-	ND				0.017	-	ND			
7440 00 0	T1 111 7 7 7 7	C-Zone	ND	0.02	0.015	1	7 0.00087 0.0009	0.0000	ND	0.02	0.015	0.00-	0.0005	0.0000
7440-28-0	Thallium [µg/L]	D-Zone	ND	ND	ND	()()()5/		0.0009	ND	ND	ND	0.005	0.005 0.00087	0.0009
			ND	ND ND		0.00007		ND	ND	ND	1			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	/ell Data**	
CASRN	Constituent [[]nite]	Lavier	Concent	ration within	Capture	Estimated C	oncentrations fo	or Production	Concent	ration within	Capture	Estimated C	oncentrations fo	or Production
CASKIN	Constituent [Units]	Layer		Zone			Wells			Zone			Wells	
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	ND				ND	-	ND			
28249-77-6	Thiobencarb [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
20245 11 0	Thioberteans [µg/L]	D-Zone	ND	ND	ND		ND	IND	ND	ND	ND	ND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	1.4	-	18				1.2	-	18			
THM	THMs, Total [µg/L]	C-Zone	3.5	11	3.3	1.2	0.55	0.59	2.6	2.6	1.9	0.99	0.17	0.49
		D-Zone	1.3	1.3	1.3		0.55	0.00	1.1	1.1	1.1	0.55	<b>3</b>	0.15
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.064	-	0.055	-			0.061	-	0.055			
7440-29-1	Thorium [µg/L]	C-Zone	ND	0.11	0.13	0.019	0.0048	0.0086	ND	0.11	0.098	0.018	0.0048	0.0067
		D-Zone	NA	NA	NA	-			NA	NA	NA			
		Sunnyside	NA 800	NA	NA 1				NA 160	NA	NA 1			
		B-Zone C-Zone	800	24	1				160 3.7	12	1			
ODOR	DDOR Threshold Odor Number [Units]	D-Zone	8	8	8	250	8.7	7.5	8	8	8	54	8.2	7.5
		Sunnyside	8	8	8	-			8	8	8			
		B-Zone	1.2	-	ND				0.51	-	ND			
		C-Zone	1.3	0.34	ND	-			0.51	0.27	ND			
108-88-3	Toluene [µg/L]	D-Zone	ND	ND	ND	0.58	0.015	ND	ND	ND	ND	0.24	0.012	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	1100000	-	990000				1100000	-	990000			
		C-Zone	1000000	1600000	890000				850000	1100000	860000			
TDS	Total Dissolved Solids [µg/L]	D-Zone	680000	680000	680000	790000	510000	500000	650000	650000	650000	730000	490000	490000
		Sunnyside	450000	450000	450000				450000	450000	450000			
		B-Zone	1200	-	580				890	-	580			
TOC	T. 10	C-Zone	1200	21000	490	550	010	20	810	11000	420	400	400	25
TOC	Total Organic Carbon [µg/L]	D-Zone	NA	NA	NA	550	910	39	NA	NA	NA	400	490	35
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
1336-36-3	Total PCBs [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
1330-30-3	rotal r Cbs [μg/L]	D-Zone	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	2400	-	91000				1300	-	91000			
TSuS	Total Suspended Solids [µg/L]	C-Zone	20000	3000	150000	4100	130	10000	12000	2500	75000	2400	110	6000
. 5 3.5	. ota: ouspoilada ooilas [µg/ -]	D-Zone	NA	NA	NA				NA	NA	NA	00		
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
1330-20-7	Total Xylenes [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	, ., 5, -	D-Zone	ND	ND	ND				ND	ND	ND			
		Sunnyside	ND	ND	ND				ND	ND	ND			+
		B-Zone	ND	- ND	ND	-			ND ND	- ND	ND ND			
8001-35-2	Toxaphene [µg/L]	C-Zone D-Zone	ND ND	ND ND	ND	ND	ND	ND	ND ND	ND ND	ND ND	ND	ND	ND
			ND ND	ND ND ND NI	D ND ND	ND ND	ND ND	ND ND		ND ND	IND			
		Sunnyside	שאו	ND	ND				ND	ND	ND			







				Us	ing UCL95 fo	or Monitoring V	Vell Data*			Us	ing Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concent	ration within Zone	Capture	Estimated Co	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	ND	-	NA				ND	-	NA			
TDUG	TDI la fina /II	C-Zone	NA	ND	NA	ND	ND	NA	NA	ND	NA	ND	ND	NIA
TPHg	TPHg [µg/L]	D-Zone	NA	NA	NA	ND	ND	INA	NA	NA	NA	ND	ND	NA
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
156-60-5	trans-1,2-Dichloroethene [µg/L]	C-Zone	ND	0.97	0.63	ND	0.042	0.037	ND	0.61	0.57	ND	0.026	0.033
130 00 3	trans 1,2 Dichloroetherie [µg/L]	D-Zone	ND	ND	ND	ND	0.042	0.037	ND	ND	ND	ND	0.020	0.033
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND			
10061-02-6	trans-1,3-Dichloropropene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
10001 02 0	tidiis 1,5 Biemoropropene [µg/ 2]	D-Zone	ND	ND	ND	110	145	110	ND	ND	ND		110	110
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND	-		
110-57-6	trans-1,4-Dichloro-2-butene [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
	tians 1,1 Diamete 2 Saterio [pg, -]	D-Zone	ND	ND	ND				ND	ND	ND	)		
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND	-	ND	<u> </u>  -		
24017-47-8	Triazofos [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
		D-Zone	ND	ND	ND				ND	ND	ND	-		IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	NA	-	NA				NA	-	NA	-		
75-96-7	Tribromoacetic acid [µg/L]	C-Zone	NA	ND	NA	ND	ND	ND	NA	ND	NA	ND	ND	ND
	3· ·	D-Zone	ND	ND	ND				ND	ND	ND	-		שאו
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	- 0.07	ND				ND	- 0.07	ND	-		
76-03-9	Trichloroacetic acid [µg/L]	C-Zone	ND	0.87	ND	ND	0.038	ND	ND	0.87	ND	ND	0.038	ND
	. •	D-Zone	ND	ND	ND				ND	ND	ND	-		
		Sunnyside	ND 2.100	ND	ND				ND	ND	ND C F 0.7			
		B-Zone	2.196	(10.24)	7.497				1.736	12.01	6.587	1		
79-01-6	Trichloroethene (TCE) [µg/L]	C-Zone D-Zone	126.9 1.5	19.24 1.5	3.838 1.5	22.2	0.92	0.44	63.6 1.5	13.01 1.5	3.104 1.5	11.4	0.65	0.38
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	0.48	-	ND				0.48	-	ND ND			
		C-Zone	0.48	0.62	ND				0.48	0.52	ND	-		
75-69-4	Trichlorofluoromethane [µg/L]	D-Zone	ND	ND	ND	0.22	0.027	ND	ND	ND	ND	0.22	0.023	ND
		Sunnyside	ND	ND	ND				ND	ND	ND	-		
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND	1		
3380-34-5	Triclosan [μg/L]	D-Zone	NA	NA NA	NA	ND	ND	ND	NA	NA	NA NA	ND	ND	ND
		Sunnyside	NA	NA	NA				NA	NA	NA NA	1		
		B-Zone	ND	-	ND				ND	-	ND			
		C-Zone	ND	ND	ND				ND	ND	ND	1		
1582-09-8	Trifluralin [μg/L]		ND	ND	ND	ND	ND	ID ND	ND	ND	ND	ND	ND	ND
		D-Zone Sunnyside	ND	ND	ND				ND	ND	ND	1		







				Us	ing UCL95 f	or Monitoring \	Well Data*			Usi	ng Mean fo	r Monitoring W	ell Data**	
CASRN	Constituent [Units]	Layer	Concent	ration within Zone	Capture	Estimated C	oncentrations fo Wells	or Production	Concentr	ation within Zone	Capture	Estimated C	oncentrations fo Wells	r Production
			SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9	SM-4	SM-8	SM-9
		B-Zone	0.00029	-	ND				0.00029	-	ND			
738-70-5	Trim oth on vine [110/1]	C-Zone	ND	0.00048	ND	9E-05	2.1E-05	ND	ND	0.00048	ND	9E-05	2.1E-05	ND
/30-70-3	Trimethoprim [µg/L]	D-Zone	NA	NA	NA	9E-05	2.1E-U5	ND	NA	NA	NA	96-05	2.1E-05	ND
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
786-19-6	Trithion [µg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
700-19-0	πιτιίοπ [μg/ε]	D-Zone	ND	ND	ND		IND	ND	ND	ND	ND	] ND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	70	-	13				70		13			
10028-17-8	Tritium [pci/L]	C-Zone	50	40	58	51	58	58	50	39	58	51	58	58
10020-17-0	muam [pc//L]	D-Zone	ND	ND	ND		36	30	ND	ND	ND	] 31	30	30
		Sunnyside	63	63	63				63	63	63			
		B-Zone	3.1	-	38				1.5	-	38			
TLIDE	TURB Turbidity [NTU]	C-Zone	31	7.9	33	6.9	1.4	3.7	16	2.1	21	3.8	1.2	3
TOND	raibiaity [ivio]	D-Zone	1.9	1.9	1.9	0.9	1.4	5.7	1.8	1.8	1.8	3.0	1.2	
		Sunnyside	1.1	1.1	1.1				1.1	1.1	1.1			
		B-Zone	10.36	-	8.8				8.333	-	8.8			
7440-61-1	Uranium Rad [pci/L]	C-Zone	9.4	3.9	14	5.3	1.3	2.1	5.733	1.574	8.9	4	1.2	1 Ω
7440-01-1	Oranium Rad [pci/L]	D-Zone	0.81	0.81	0.81		1.5	2.1	0.81	0.81	0.81	4	1.2	1.8
		Sunnyside	1.2	1.2	1.2				1.2	1.2	1.2			
		B-Zone	3.185	-	11				2.233	-	11			
7440-62-2	Vanadium, Total [µg/L]	C-Zone	8.4	3.1	6.6	2.4	0.13	0.58	5.308	1.635	5.35	1.6	0.071	0.51
7440 02 2	vanadidiii, i Otai [μg/ L]	D-Zone	NA	NA	NA	2,4	0.13	0.50	NA	NA	NA	1.0	0.071	0.51
		Sunnyside	NA	NA	NA				NA	NA	NA			
		B-Zone	ND	-	ND				ND	-	ND			
108-05-4	Vinyl Acetate [μg/L]	C-Zone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
100-03-4	Villyi Acetate [µg/L]	D-Zone	ND	ND	ND	IND	IND	ND	ND	ND	ND	IND	ND	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	ND	-	ND				ND		ND			
75-01-4	Vinyl Chloride [µg/L]	C-Zone	ND	1.9	ND	ND	0.081	ND	ND	1.3	ND	ND	0.058	ND
7501-4	Villyi Ciliolide [µg/L]	D-Zone	ND	ND	ND	IND	0.001	IND	ND	ND	ND	ND	0.030	IND
		Sunnyside	ND	ND	ND				ND	ND	ND			
		B-Zone	6.3	-	1				3.5	-	1			
7440-66-6	Zinc [μg/L]	C-Zone	1.8	1.1	2	5.1	5.1 2.6	26 26	1.4	1	1.5	3.9	2.5	2.5
7770 00-0	Zinc [µg/L]	D-Zone	12	12	12			2.6	10	10	10	3.9	3.9 2.5	
		Sunnyside	2.1	2.1	2.1				2.1	2.1	2.1			

## Notes:

CASRN = Chemical Abstracts Service Registry Number (as applicable); μg/L = micrograms per liter; ng/L = nanograms per liter; mpN/100mL = most probable number per 100 millilitres; NTU = nephelometric turbidity units; 'ND' = non-detect; 'NA' = not available; '-' = not applicable.

\*UCL95 (95 percent upper confidence limit of the population mean) based on maximum suggested UCL95 statistics produced by ProUCL software (EPA 2015) calculated for censored data sets (using all detects and non-detects).

For the D-Zone, monitoring well SMB-1-B#1 was used for all production well capture zones; for the Sunnyside, monitoring well SMB-1 #4 was used for all production well capture zones.

For constituents (metals) where both total and dissolved fractions were reported for the same sample, values for the dissolved fraction were used.

<sup>\*\*</sup>Mean calculated for censored data sets (using all detects and non-detects) via Kaplan Meier Method (using ProUCL software; EPA 2015).

In the event either the mean or UCL95 statistics could not be calculated, the maximum value was adopted as a conservative approach.

For non-detect (ND) results a value of 0 was assumed; where data was not available (NA) for a constituent in any given layer and/or capture zone, a value of 0 was assumed.







# Appendix F Treatment Plant Influent Concentration Estimates







			Well Data*	U	sing Maximum Va	alue for 2020 Prod	duction Well Data**		
CASRN	Constituent [Units]	Estimated Con	centrations for Pr		Plant Influent Concentration		centrations For P		Plant Influent Concentration
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates
3567-62-2	1-(3,4-Dichlorophenyl)-3-methylurea [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
2327-02-8	1-(3,4-Dichlorophenyl)urea [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
630-20-6	1,1,1,2-Tetrachloroethane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
71-55-6	1,1,1-Trichloroethane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
79-34-5	1,1,2,2-Tetrachloroethane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
79-00-5	1,1,2-Trichloroethane [µg/L]	0.27	ND	ND	0.094	ND	ND	ND	ND
75-34-3	1,1-Dichloroethane (1,1-DCA) [µg/L]	0.27	0.039	0.01	0.11	0.16	0.12	ND	0.098
75-35-4	1,1-Dichloroethene (1,1-DCE) [µg/L]	1.1	0.2	0.079	0.47	1.2	0.72	ND	0.67
563-58-6	1,1-Dichloropropene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
57-14-7	1,1-Dimethylhydrazine [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
87-61-6	1,2,3-Trichlorobenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND ND
96-18-4	1,2,3-Trichloropropane (1,2,3-TCP) [µg/L]	0.037	0.015	0.014	0.022	0.0026	ND	ND	0.00091
120-82-1	1,2,4-Trichlorobenzene [µg/L]	ND	ND	ND ND	ND	ND	ND	ND	ND
95-63-6	1,2,4-Trimethylbenzene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
96-12-8	1,2-Dibromo-3-chloropropane [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
106-93-4	1,2-Dibromoethane [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
95-50-1	1,2-Dichlorobenzene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
107-06-2	1,2-Dichloroethane [µg/L]	0.083	0.034	0.0085	0.043	ND	ND	ND	ND ND
78-87-5	1,2-Dichloropropane [µg/L]	0.048	ND	ND	0.017	ND	ND	ND	ND ND
122-66-7	1,2-Diphenylhydrazine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND ND
108-67-8	1,3,5-Trimethylbenzene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
99-35-4	1,3,5-Trinitrobenzene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
541-73-1	1,3-Dichlorobenzene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
142-28-9	1,3-Dichloropropane [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
542-75-6	1,3-Dichloropropene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
99-65-0	1,3-Dinitrobenzene [µg/L]	0.044	ND	ND	0.015	ND	ND	0.32	0.096
106-46-7	1,4-Dichlorobenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
123-91-1	1,4-Dioxane [µg/L]	35.4	2.1	2.2	13.8	20	5.4	3.3	9.9
763051-92-9	11CI-PF3OUdS [µg/L]	ND	ND	ND	ND	ND ND	ND	ND	ND
57-91-0	17-a-Estradiol [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
57-63-6	17-a-Ethynylestradiol [μg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
50-28-2	17-b-Estradiol [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
124-19-6	1-Nonanal [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
934-34-9	2(3H)-Benzothiazolone [µg/L]	NA NA	0.061	NA NA	0.021	ND	ND	ND	NA NA
594-20-7	2,2-Dichloropropane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND ND
51207-31-9	2,3,7,8-Tetrachlorodibenzofuran [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
93-76-5	2,4,5-Τ [μg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
93-72-1	2,4,5-TP (Silvex) [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
95-95-4	2,4,5-Trichlorophenol [µg/L]	ND	ND	ND	ND ND	ND ND	ND	ND	ND ND
88-06-2	2,4,6-Trichlorophenol [µg/L]	ND	ND	ND	ND ND	ND ND	ND	ND	ND ND
118-96-7	2,4,6-Trinitrotoluene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
94-75-7	2,4-D [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
94-82-6	2,4-DB [μg/L]	ND	ND	ND	ND ND	ND ND	ND	ND	ND ND
120-83-2	2,4-Dichlorophenol [μg/L]	ND	ND	ND	ND ND	ND ND	ND	ND	ND
105-67-9	2,4-Dictriorophenol [µg/L]	ND	ND ND	ND	ND ND	ND ND	ND ND	ND	ND ND







CASRN	Constituent [Units]	_			Well Data*		duction Well Data**		
		Estimated Con	centrations for Pr		Plant Influent Concentration		centrations For P		Plant Influent Concentration
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates
51-28-5	2,4-Dinitrophenol [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
121-14-2	2,4-Dinitrotoluene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
606-20-2	2,6-Dinitrotoluene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
35572-78-2	2-Amino-4,6-Dinitrotoluene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
78-93-3	2-Butanone [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
110-75-8	2-Chloroethyl Vinyl Ether [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
91-58-7	2-Chloronaphthalene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
95-57-8	2-Chlorophenol [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
95-49-8	2-Chlorotoluene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
591-78-6	2-Hexanone [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
91-57-6	2-Methylnaphthalene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
95-48-7	2-Methylphenol [μg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
88-74-4	2-Nitroaniline [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
88-75-5	2-Nitrophenol [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
88-72-2	2-Nitrotoluene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
7642-04-8	2-Octene, (Z)- [µg/L]	9	NA	6	5	ND	ND	34	10
84989-04-8	3 & 4-Methylphenol [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
91-94-1	3,3'-Dichlorobenzidine [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
95-76-1	3,4-Dichloroaniline [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
51-36-5	3,5-Dichlorobenzoic acid [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
16655-82-6	3-Hydroxycarbofuran [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
99-09-2	3-Nitroaniline [μg/L]	ND	ND	ND	ND ND	ND	ND	ND ND	ND ND
99-08-1	3-Nitrotoluene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND ND	ND ND
72-54-8	4,4´-DDD [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
72-55-9	4,4´-DDE [µg/L]	ND	ND	ND	ND ND	ND	ND ND	ND	ND
50-29-3	4,4´-DDT [µg/L]	ND	ND	ND	ND ND	ND	ND	ND ND	ND ND
534-52-1	4,6-Dinitro-2-Methyl phenol [µg/L]	ND	ND	ND	ND ND	ND	ND ND	ND	ND
19406-51-0	4-Amino-2,6-Dinitrotoluene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
101-55-3	4-Bromophenyl Phenyl Ether [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
59-50-7	4-Chloro-3-Methylphenol [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
106-47-8	4-Chloroaniline [µg/L]	ND	ND	ND	ND ND	ND	ND ND	ND	ND ND
7005-72-3	4-Chlorophenyl Phenyl Ether [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
106-43-4	4-Chlorotoluene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
108-10-1	4-Methyl-2-pentanone [μg/L]	ND	ND	ND	ND ND	ND	ND ND	ND	ND ND
100-01-6	4-Nitroaniline [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
100-02-7	4-Nitrophenol [µg/L]	ND	ND	ND	ND ND	ND	ND ND	ND	ND ND
99-99-0	4-Nitrotoluene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
104-40-5	4-Nonylphenol [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
1806-26-4	4-Octylphenol [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
140-66-9	4-tert-Octylphenol [µg/L]	ND	0.00061	ND	0.00021	ND	ND	ND	ND
2315-61-9	4-tert-Octylphenol diethoxylate [µg/L]	ND	0.00001 ND	ND	ND	ND	ND ND	ND ND	ND
1173019-48-1	4-tert-Octylphenol monoethoxylate [µg/L]	0.0027	0.0031	ND	0.002	ND	ND ND	ND	ND
756426-58-1	9CI-PF3ONS [µg/L]	ND	ND	ND ND	0.002 ND	ND ND	ND ND	ND	ND ND
83-32-9	Acenaphthene [µg/L]	ND	ND	ND	ND	ND	ND	ND ND	ND
208-96-8	Acenaphthene [µg/L] Acenaphthylene [µg/L]	ND	ND ND	ND ND	ND ND	ND ND	ND ND	ND ND	ND







			Using UCL9	5 for Monitoring	Well Data*	U	sing Maximum Va	alue for 2020 Pro	duction Well Data**
CASRN	Constituent [Units]	Estimated Con	centrations for Pr	oduction Wells	Plant Influent Concentration	Observed Con	centrations For P	Production Well	Plant Influent Concentration
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates
75-07-0	Acetaldehyde [µg/L]	0.21	ND	ND	0.072	ND	ND	ND	ND
103-90-2	Acetaminophen [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
34256-82-1	Acetochlor [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
187022-11-3	Acetochlor ESA [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
194992-44-4	Acetochlor OA [µg/L]	ND	ND	ND	ND	0.0068	ND	ND	0.0024
67-64-1	Acetone [µg/L]	15	7.3	6.8	9.8	790	ND	22	280
75-05-8	Acetonitrile [µg/L]	0.29	0.22	0.21	0.24	7.9	ND	ND	2.8
50594-66-6	Acifluorfen [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
107-02-8	Acrolein [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
107-13-1	Acrylonitrile [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
958445-44-8	ADONA [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
15972-60-8	Alachlor [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
142363-53-9	Alachlor ESA [µg/L]	ND	0.00091	ND	0.00032	ND	ND	ND	ND
171262-17-2	Alachlor OA [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
116-06-3	Aldicarb [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1646-88-4	Aldicarb Sulfone [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1646-87-3	Aldicarb sulfoxide [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
309-00-2	Aldrin [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
ALK	Alkalinity (as CaCO3) [µg/L]	310000	230000	240000	260000	360000	250000	430000	340000
ALKC-1	Alkalinity, Carbonate as CaCO3 [µg/L]	ND	2700	ND	950	ND	ND	ND	ND
ALKH-1	Alkalinity, Hydroxide as CaCO3 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
107-05-1	Allyl Chloride [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
319-84-6	alpha-BHC [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
5103-71-9	alpha-Chlordane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7429-90-5	Aluminum, Total [µg/L]	15.4	5.2	51.3	22.6	3.4	1.9	2.5	2.6
7664-41-7	Ammonia [µg/L]	35	52	ND	30	ND	34	ND	12
26787-78-0	Amoxicillin [μg/L]	ND	ND	ND	ND	ND	0.0055	ND	0.0019
62-53-3	Aniline [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
120-12-7	Anthracene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7440-36-0	Antimony [µg/L]	0.15	0.21	0.21	0.19	0.072	0.093	0.15	0.1
12674-11-2	Aroclor 1016 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
11104-28-2	Aroclor 1221 [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
11141-16-5	Aroclor 1232 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
53469-21-9	Aroclor 1242 [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
12672-29-6	Aroclor 1248 [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
11097-69-1	Aroclor 1254 [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
11096-82-5	Aroclor 1260 [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
22541-54-4	Arsenic III [µg/L]	1.9	4.3	4.1	3.4	ND	0.17	ND	0.06
17428-41-0	Arsenic V [μg/L]	0.51	0.032	0.034	0.2	0.99	0.49	0.98	0.81
7440-38-2	Arsenic, Total [μg/L]	0.58	1.64	0.16	0.83	0.71	0.69	0.74	0.71
1332-21-4	Asbestos [MFL]	ND	ND	ND	ND	ND	ND	ND	ND
29122-68-7	Atenolol [μg/L]	0.002	0.000014	ND	0.0007	ND	0.00028	ND	0.000098
134523-00-5	Atorvastatin [µg/L]	0.00013	ND	ND	0.000045	0.0011	ND	ND	0.00039
1912-24-9	Atrazine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
83905-01-5	Azithromycin [µg/L]	0.0044	ND	ND	0.0015	0.0057	ND	ND	0.002







			Using UCL9	5 for Monitoring	Well Data*	U	sing Maximum Va	alue for 2020 Prod	duction Well Data**
CASRN	Constituent [Units]	Estimated Con	centrations for Pr		Plant Influent Concentration		centrations For P		Plant Influent Concentration
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates
103-33-3	Azobenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7440-39-3	Barium, Dissolved [µg/L]	19.8	0.82	4	8.4	45	60	34	47
7440-39-3	Barium, Total [µg/L]	20.4	1.26	4.3	8.9	60	58	34	52
25057-89-0	Bentazon [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
100-52-7	Benzaldehyde [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
71-43-2	Benzene [µg/L]	0.039	0.068	ND	0.037	ND	ND	ND	ND
92-87-5	Benzidine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
56-55-3	Benzo[a]anthracene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
50-32-8	Benzo[a]pyrene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
205-99-2	Benzo[b]fluoranthene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
191-24-2	Benzo[g,h,i]perylene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
207-08-9	Benzo[k]fluoranthene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
65-85-0	Benzoic Acid [µg/L]	ND	0.29	ND	0.1	ND	ND	ND	ND
95-16-9	Benzothiazole [µg/L]	NA	0.37	NA	0.13	ND	ND	ND	NA NA
100-51-6	Benzyl Alcohol [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
85-68-7	Benzyl butyl phthalate [µg/L]	0.51	0.17	0.29	0.32	3.7	ND	ND	1.3
7440-41-7	Beryllium [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
319-85-7	beta-BHC [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
71-52-3	Bicarbonate Alkalinity as HCO3 [µg/L]	390000	280000	290000	320000	440000	300000	520000	420000
BOD	Biochemical Oxygen Demand [µg/L]	ND	300	ND	100	ND	ND	ND	ND
111-91-1	bis(2-chloroethoxy)methane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
111-44-4	bis(2-chloroethyl)ether [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
108-60-1	bis(2-chloroisopropyl)ether [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
103-23-1	Bis(2-ethylhexyl)adipate [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
117-81-7	bis(2-ethylhexyl)phthalate [µg/L]	0.38	0.7	0.63	0.57	ND	ND	ND	ND
80-05-7	Bisphenol A [µg/L]	0.0029	0.0056	0.00014	0.003	0.041	0.0027	0.0014	0.016
7440-42-8	Boron, Total [µg/L]	63.3	26	11.4	34.7	130	140	130	130
314-40-9	Bromacil [μg/L]	0.062	ND	ND	0.022	ND	ND	ND	ND
15541-45-4	Bromate [µg/L]	10	7.1	0.76	6.3	ND	ND	ND	ND
24959-67-9	Bromide [µg/L]	330	250	31	210	540	230	850	520
108-86-1	Bromobenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
5589-96-8	Bromochloroacetic acid [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
74-97-5	Bromochloromethane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
71133-14-7	Bromodichloroacetic acid [µg/L]	ND	ND	ND	ND	ND	ND	ND	NA
75-27-4	Bromodichloromethane [µg/L]	0.077	ND	ND	0.027	0.095	ND	0.1	0.063
75-25-2	Bromoform [µg/L]	ND	ND	ND	ND	1.5	ND	ND	0.53
74-83-9	Bromomethane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
23184-66-9	Butachlor [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
123-72-8	Butanal [µg/L]	0.82	ND	ND	0.29	ND	ND	ND	ND
7440-43-9	Cadmium [µg/L]	0.054	0.004	0.0041	0.021	0.083	ND	0.12	0.065
58-08-2	Caffeine [µg/L]	0.011	0.0026	0.00025	0.0048	0.001	0.00091	0.031	0.01
7440-70-2	Calcium [µg/L]	120000	73000	76000	90000	140000	91000	180000	140000
133-06-2	Captan [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
298-46-4	Carbamazepine [µg/L]	ND	ND	ND	ND	ND	ND	0.00009	0.000027
63-25-2	Carbaryl [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND







			Usina UCL9	5 for Monitoring	Well Data*	U	sing Maximum Va	alue for 2020 Proc	luction Well Data**
CASRN	Constituent [Units]	Estimated Con	centrations for Pr		Plant Influent Concentration		centrations For P		Plant Influent Concentration
3.131	, , , , , , , , , , , , , , , , , , , ,	SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates
86-74-8	Carbazole [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1563-66-2	Carbofuran [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
75-15-0	Carbon Disulfide [µg/L]	ND	0.011	ND	0.0039	ND	ND	ND	ND
56-23-5	Carbon Tetrachloride [µg/L]	0.36	0.044	0.024	0.15	0.13	0.15	ND	0.098
133-90-4	Chloramben [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7775-09-9	Chlorate [µg/L]	54.3	1.4	11.1	22.8	91	ND	93	60
57-74-9	Chlordane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
16887-00-6	Chloride [µg/L]	93000	86000	65000	82000	120000	110000	130000	120000
76-13-1	Chlorinated Fluorocarbon (Freon 113) [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
7782-50-5RF	Chlorine Residual, Free [µg/L]	4.3	0.56	5.7	3.4	35	9	2	16
TOTAL-CHLORINE	Chlorine Residual, Total [µg/L]	11	0.95	8.7	6.8	ND	ND	ND	NA NA
14998-27-7	Chlorite [µg/L]	1.6	3.5	2.7	2.6	13	14	ND	9.5
79-11-8	Chloroacetic Acid [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
108-90-7	Chlorobenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
5278-95-5	Chlorodibromoacetic acid [µg/L]	ND	ND	ND	ND	ND	ND	ND	NA NA
75-00-3	Chloroethane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
67-66-3	Chloroform [µg/L]	1.1	0.18	0.47	0.59	2.2	0.66	0.9	1.3
74-87-3	Chloromethane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1897-45-6	Chlorothalonil [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
101-21-3	Chlorpropham [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND ND
18540-29-9	Chromium, Hexavalent [µg/L]	0.46	0.017	0.39	0.28	1.1	0.81	0.25	0.74
7440-47-3	Chromium, Total [µg/L]	51.5	0.056	14.4	22.4	0.93	0.051	0.24	0.42
218-01-9	Chrysene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
85721-33-1	Ciprofloxacin [µg/L]	0.0042	0.00012	0.00084	0.0018	0.11	ND	0.0065	0.04
156-59-2	cis-1,2-Dichloroethene (cis-1,2-DCE) [µg/L]	0.22	2.1	0.047	0.84	0.45	0.22	ND	0.23
10061-01-5	cis-1,3-Dichloropropene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
5103-73-1	cis-Nonachlor [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7440-48-4	Cobalt [µg/L]	2.7	0.004	0.077	0.98	0.09	0.16	0.071	0.11
CLR	Color [Units]	ND	ND	ND	ND	ND	ND	ND	ND
CLR	Color [Units]	ND	ND	ND	ND	ND	ND	ND	ND
7440-50-8	Copper [µg/L]	0.84	0.22	0.3	0.46	3.3	0.14	2.3	1.9
486-56-6	Cotinine [µg/L]	0.0015	0.0012	0.00016	0.00099	ND	ND	0.0012	0.00036
123-73-9	Crotonaldehyde [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
21725-46-2	Cyanazine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
57-12-5	Cyanide [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1122-82-3	Cyclohexane, isothiocyanato- [µg/L]	NA	0.043	NA	0.015	ND	ND	ND	NA
108-94-1	Cyclohexanone [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
75-99-0	Dalapon [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1861-32-1	DCPA [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
112-31-2	Decanal [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
134-62-3	DEET [µg/L]	0.0011	0.00016	0.00019	0.0005	0.0043	0.001	0.0008	0.0021
319-86-8	delta-BHC [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
439-14-5	Diazepam [μg/L]	0.000063	ND	ND	0.000022	ND	ND	ND	ND
333-41-5	Diazinon [µg/L]	ND	0.0024	ND	0.00083	ND	ND	ND	ND
53-70-3	Dibenzo[a,h]anthracene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND







			Using UCLS	5 for Monitoring	Well Data*	U	sing Maximum V	alue for 2020 Pro	duction Well Data**
CASRN	Constituent [Units]	Estimated Con	centrations for Pr		Plant Influent Concentration		centrations For P		Plant Influent Concentration
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates
132-64-9	Dibenzofuran [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
631-64-1	Dibromoacetic acid [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
124-48-1	Dibromochloromethane [µg/L]	ND	3	ND	1.1	ND	ND	ND	ND
74-95-3	Dibromomethane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1918-00-9	Dicamba [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
3400-09-7	Dichloramine [μg/L]	2.4	0.17	0.88	1.2	7	4	2	4.5
79-43-6	Dichloroacetic Acid [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
75-71-8	Dichlorodifluoromethane [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
120-36-5	Dichloroprop [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
15307-86-5	Diclofenac [μg/L]	0.00025	ND	ND	0.000088	ND	ND	ND	ND
60-57-1	Dieldrin [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
84-66-2	Diethyl phthalate [μg/L]	0.14	0.19	0.16	0.16	ND	ND	ND	ND
56-53-1	Diethylstilbestrol [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
108-20-3	Di-isopropyl ether [µg/L]	ND	0.053	ND	0.019	ND	ND	ND	ND
60-51-5	Dimethoate [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
131-11-3	Dimethyl phthalate [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
84-74-2	Di-n-butyl-phthalate [μg/L]	0.26	0.25	0.29	0.27	0.23	ND	0.19	0.14
117-84-0	Di-n-octyl phthalate [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
88-85-7	Dinoseb [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
957-51-7	Diphenamid [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
85-00-7	Diquat [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
298-04-4	Disulfoton [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
330-54-1	Diuron [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
959-98-8	Endosulfan I [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
33213-65-9	Endosulfan II [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1031-07-8	Endosulfan sulfate [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
145-73-3	Endothall [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
72-20-8	Endrin [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7421-93-4	Endrin Aldehyde [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
53494-70-5	Endrin Ketone [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
481-30-1	Epitestosterone [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
759-94-4	EPTC [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
50-27-1	Estriol [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
53-16-7	Estrone [µg/L]	0.000084	ND	0.000045	0.000043	ND	ND	ND	ND
2991-50-6	EtFOSAA [μg/L]	0.00027	ND	ND	0.000094	ND	ND	ND	ND
64-17-5	Ethanol [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
563-12-2	Ethion [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
97-63-2	Ethyl Methacrylate [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
637-92-3	Ethyl tert-butyl ether [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
100-41-4	Ethylbenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
107-21-1	Ethylene Glycol [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
206-44-0	Fluoranthene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
86-73-7	Fluorene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
16984-48-8	Fluoride, Total [mg/L]	0.15	0.01	0.024	0.06	0.29	0.24	0.33	0.28
54910-89-3	Fluoxetine [µg/L]	0.00044	ND	0.000012	0.00016	0.00094	ND	0.0011	0.00066







			Using UCL9	5 for Monitoring	Well Data*	U	sing Maximum Va	alue for 2020 Prod	duction Well Data**
CASRN	Constituent [Units]	Estimated Con	centrations for Pr		Plant Influent Concentration		centrations For P		Plant Influent Concentration
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates
944-22-9	Fonofos [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
50-00-0	Formaldehyde [µg/L]	ND	0.039	ND	0.013	ND	ND	ND	ND
1222-05-5	Galaxolide (HHCB) [μg/L]	0.021	0.001	0.0013	0.0081	0.16	0.0048	0.14	0.1
58-89-9	gamma-BHC [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
5566-34-7	gamma-Chlordane [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
8032-32-4	Gasoline Range Organics [µg/L]	180	4.3	8.8	67	160	ND	140	98
25812-30-0	Gemfibrozil [µg/L]	0.000033	0.0000052	ND	0.000013	ND	0.000098	ND	0.000034
107-22-2	Glyoxal [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1071-83-6	Glyphosate [µg/L]	0.38	0.21	0.12	0.24	ND	ND	ND	ND
12587-46-1	Gross Alpha [pci/L]	3.2	0.09	0.56	1.3	3.5	4.4	9.6	5.6
12587-47-2	Gross Beta [pci/L]	2.6	0.36	0.83	1.3	8.1	4.4	6.7	6.4
HAA5	HAA5, Total [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
Hardt-1	Hardness, total as CaCO3 [µg/L]	510000	290000	300000	370000	640000	390000	830000	610000
76-44-8	Heptachlor [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
1024-57-3	Heptachlor Epoxide [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
111-71-7	Heptanal [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
118-74-1	Hexachlorobenzene [µg/L]	0.00023	0.0039	0.00025	0.0015	ND	ND	ND	ND
87-68-3	Hexachlorobutadiene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
77-47-4	Hexachlorocyclopentadiene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
67-72-1	Hexachloroethane [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
66-25-1	Hexanal [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
13252-13-6	HFPO-DA [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
2691-41-0	HMX [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
302-01-2	Hydrazine [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
15687-27-1	Ibuprofen [μg/L]	0.0018	0.00021	0.000039	0.00072	ND	ND	0.0013	0.00039
193-39-5	Indeno[1,2,3-cd]pyrene [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
20461-54-5	lodide [μg/L]	7.7	3.3	0.49	4	6.9	34	1.4	15
74-88-4	lodomethane [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
73334-07-3	lopromide [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7439-89-6	lron [μg/L]	7.2	120	8.7	48	1.4	ND	ND	0.49
78-59-1	Isophorone [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
67-63-0	Isopropyl Alcohol [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
98-82-8	Isopropylbenzene [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7439-92-1	Lead, Total [µg/L]	0.043	0.016	0.03	0.03	0.22	0.13	0.2	0.18
330-55-2	Linuron [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7439-93-2	Lithium [µg/L]	ND	1.1	ND	0.38	ND	ND	ND	ND
179601-23-1	m,p-Xylene [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7439-95-4	Magnesium [μg/L]	58000	26000	27000	38000	67000	42000	91000	65000
7439-96-5	Manganese, Total [μg/L]	107.2	4.8	6	41	33	46	3.5	29
2355-31-9	MeFOSAA [μg/L]	0.00022	ND	ND	0.00076	ND	ND	ND	ND
57-53-4	Meprobamate [µg/L]	0.00037	ND	ND	0.00013	ND	ND	ND	ND .
7439-97-6	Mercury [μg/L]	0.017	0.019	0.018	0.018	ND	0.023	ND	0.0081
126-98-7	Methacrylonitrile [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
76-99-3	Methadone [µg/L]	0.00011	0.0000027	ND	0.00039	0.00016	ND	ND	0.000056
67-56-1	Methanol [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND







			Using UCL9	5 for Monitoring	Well Data*	U	sing Maximum Va	alue for 2020 Prod	luction Well Data**
CASRN	Constituent [Units]	Estimated Con	centrations for Pr		Plant Influent Concentration		centrations For P		Plant Influent Concentration
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates
2032-65-7	Methiocarb [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
16752-77-5	Methomyl [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
72-43-5	Methoxychlor [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
78-98-8	Methyl Glyoxal [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
80-62-6	Methyl Methacrylate [µg/L]	0.039	ND	0.0015	0.014	ND	ND	ND	ND
1634-04-4	Methyl tert-butyl ether (MTBE) [µg/L]	0.16	0.17	ND	0.11	ND	ND	ND	ND
MBAS	Methylene blue active substances (MBAS) [µg/L]	52	79	76	69	ND	49	ND	17
75-09-2	Methylene Chloride [μg/L]	0.4	0.72	0.66	0.59	ND	ND	ND	ND
51218-45-2	Metolachlor [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
171118-09-5	Metolachlor ESA [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
152019-73-3	Metolachlor OA [µg/L]	ND	ND	ND	ND	0.0039	ND	ND	0.0014
21087-64-9	Metribuzin [µg/L]	ND	0.0034	ND	0.0012	ND	ND	ND	ND
2385-85-5	Mirex [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
2212-67-1	Molinate [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND ND
7439-98-7	Molybdenum [µg/L]	14	2	1.2	6	7.4	5.5	8.3	7
79-08-3	Monobromoacetic acid [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
10599-90-3	Monochloramine [µg/L]	3.8	0.35	2.4	2.2	6	6	4	5.4
60-34-4	Monomethylhydrazine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
57-27-2	Morphine [μg/L]	0.00063	0.000078	0.000072	0.00027	ND	ND	0.00058	0.00017
91-20-3	Naphthalene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND ND
22204-53-1	Naproxen [μg/L]	0.00087	ND	ND	0.00031	ND	ND	0.0011	0.00033
104-51-8	n-Butylbenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
7440-02-0	Nickel, Total [µg/L]	74.3	0.16	11.3	29.4	0.54	1.7	0.9	1.1
14797-55-8	Nitrate as N [mg/L]	5.6	0.23	0.6	2.2	5.8	0.62	3.3	3.2
14797-65-0	Nitrite as N [μg/L]	51.7	2.8	6.2	20.9	ND	ND	ND	ND
98-95-3	Nitrobenzene [µg/L]	ND	ND	ND	ND ND	ND	ND	ND	ND
55-18-5	N-Nitrosodiethylamine (NDEA)	0.2	0.061	0.062	0.11	0.0015	ND	ND	0.00053
62-75-9	N-Nitrosodimethylamine [µg/L]	0.0003	ND	ND	0.00011	ND	ND	ND	ND
924-16-3	n-Nitrosodi-n-butylamine [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND
621-64-7	n-Nitrosodi-n-propylamine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
86-30-6	n-Nitrosodiphenylamine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
10595-95-6	N-Nitrosomethylethylamine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
59-89-2	n-Nitrosomorpholine [μg/L]	0.012	0.00014	0.00012	0.0041	0.012	ND	0.039	0.016
100-75-4	n-Nitrosopiperidine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
930-55-2	n-Nitrosopyrrolidine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
NN	NO2+NO3 as N [mg/L]	6.9	1	1.4	3.2	5.8	0.62	3.3	3.2
25154-52-3	Nonylphenol [μg/L]	0.12	0.19	ND	0.11	0.016	ND	ND	0.0056
20427-84-3	Nonylphenol diethoxylate [μg/L]	0.014	0.14	ND	0.054	0.05	ND	ND	0.018
27986-36-3	Nonylphenol monoethoxylate [µg/L]	0.022	0.1	0.00076	0.043	0.011	ND	ND	0.0039
103-65-1	n-Propylbenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND
136777-61-2	o&p-Xylene [μg/L]	ND	ND	ND	ND	ND	ND	ND	NA
111-66-0	Octene-1 [µg/L]	4.2	NA	NA	1.5	ND	ND	ND	NA
OILGREASE	Oil & Grease (HEM) [µg/L]	450	ND	ND	160	3100	2200	ND	1900
14265-44-2	o-Phosphate as P [µg/L]	37	4.1	6.7	17	80	150	98	110
23135-22-0	Oxamyl [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND







			Using UCL9	5 for Monitoring	Well Data*	Using Maximum Value for 2020 Production Well Data**					
CASRN	Constituent [Units]	Estimated Con	centrations for Pr	oduction Wells	Plant Influent Concentration	Observed Con	centrations For P	roduction Well	Plant Influent Concentration		
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates		
131-57-7	Oxybenzone [µg/L]	0.0042	0.0013	0.00066	0.0021	0.032	0.0039	0.0041	0.014		
95-47-6	o-Xylene [μg/L]	ND	0.013	ND	0.0047	ND	ND	ND	ND		
4685-14-7	Paraquat [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
189084-64-8	PBDE-100 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
182677-30-1	PBDE-138 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
68631-49-2	PBDE-153 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
207122-15-4	PBDE-154 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
147217-75-2	PBDE-17 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
41318-75-6	PBDE-28 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
5436-43-1	PBDE-47 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
243982-82-3	PBDE-49 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
182346-21-0	PBDE-85 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
60348-60-9	PBDE-99 [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
76-01-7	Pentachloroethane [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
82-68-8	Pentachloronitrobenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
87-86-5	Pentachlorophenol [μg/L]	ND	0.019	ND	0.0065	ND	ND	ND	ND		
110-62-3	Pentanal [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
14797-73-0	Perchlorate [µg/L]	2.6	0.087	0.18	0.98	1.6	ND	ND	0.56		
375-73-5	PFBS [μg/L]	0.0022	0.000061	0.00017	0.00085	0.0028	ND	0.0051	0.0025		
335-76-2	PFDA [µg/L]	0.0019	ND	0.00012	0.00069	ND	ND	ND	ND		
307-55-1	PFDoA [μg/L]	0.0003	0.000078	0.000033	0.00014	ND	ND	ND	ND		
375-85-9	PFHpA [µg/L]	0.0022	0.00039	0.0004	0.001	0.0018	0.0018	0.0027	0.0021		
307-24-4	PFHxA [μg/L]	0.0047	0.00015	0.00026	0.0018	0.0049	ND	0.012	0.0053		
355-46-4	PFHxS [μg/L]	0.0024	0.000056	0.00021	0.00093	0.003	ND	0.0048	0.0025		
375-95-1	PFNA [µg/L]	0.0013	ND	ND	0.00044	ND	ND	0.00062	0.00019		
335-67-1	PFOA [ng/L]	1.1	0.039	0.12	0.44	0.96	ND	2.9	1.2		
1763-23-1	PFOS [ng/L]	ND	ND	0.075	0.022	ND	ND	ND	ND		
376-06-7	PFTeDA [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
72629-94-8	PFTrDA [μg/L]	ND	0.000018	ND	0.0000062	ND	ND	ND	ND		
2058-94-8	PFUnA [μg/L]	0.0013	0.000034	0.00018	0.00051	ND	ND	ND	ND		
TPHd	PHC AS DIESEL FUEL [μg/L]	NA	4.3	NA	1.5	ND	ND	ND	NA		
85-01-8	Phenanthrene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
108-95-2	Phenol [μg/L]	ND	0.052	ND	0.018	ND	ND	ND	ND		
57-41-0	Phenytoin (Dilantin) [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
7723-14-0	Phosphorus [μg/L]	200	13	55	90	620	460	260	460		
85-44-9	Phthalic Anhydride [μg/L]	3.3	0.65	NA	1.4	ND	ND	ND	NA		
1918-02-1	Picloram [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
99-87-6	p-lsopropyltoluene [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
7440-09-7	Potassium [μg/L]	4600	5200	4900	4900	3100	2500	3000	2900		
55268-74-1	Praziquantel [μg/L]	ND	0.00029	ND	0.0001	ND	ND	ND	ND		
125-33-7	Primidone [μg/L]	0.00087	ND	0.000032	0.00031	0.0012	ND	0.0034	0.0014		
57-83-0	Progesterone [μg/L]	ND	0.000014	ND	0.000005	ND	ND	ND	ND		
1610-18-0	Prometon [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
7287-19-6	Prometryn [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
1918-16-7	Propachlor (2-Chloro-n-(1-Methylethy)-n-phenylac [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		







			Using UCL9	5 for Monitoring	Well Data*	Using Maximum Value for 2020 Production Well Data**					
CASRN	Constituent [Units]	Estimated Con	centrations for Pr		Plant Influent Concentration		centrations For P	roduction Well Plant Influent Concentration			
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates		
123-38-6	Propanal [μg/L]	0.6	ND	ND	0.21	ND	ND	ND	ND		
114-26-1	Propoxur (Baygon) [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
129-00-0	Pyrene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
110-86-1	Pyridine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
91-22-5	Quinoline [µg/L]	0.0012	0.00026	0.000088	0.00054	ND	ND	ND	ND		
13982-63-3	Radium 226 [pci/L]	0.15	0.02	0.014	0.063	0.64	0.06	ND	0.25		
15262-20-1	Radium 228 [pci/L]	0.33	0.043	0.021	0.14	0.24	0.45	0.68	0.45		
121-82-4	RDX [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
69-72-7	Salicylic Acid [µg/L]	0.27	0.032	0.029	0.12	0.013	1	0.85	0.61		
135-98-8	sec-Butylbenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
7782-49-2	Selenium, Total [µg/L]	2.1	0.031	0.28	0.85	4.6	0.51	4.4	3.1		
7631-86-9	Silica [µg/L]	16000	1400	2600	6900	36000	46000	32000	38000		
7440-22-4	Silver [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
122-34-9	Simazine [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
7440-23-5	Sodium [μg/L]	78000	52000	53000	61000	82000	72000	74000	76000		
10098-97-2	Strontium-90 [pci/L]	0.25	0.27	0.25	0.255	0.61	0.84	0.26	0.58		
100-42-5	Styrene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
56038-13-2	Sucralose [µg/L]	0.0014	ND	ND	0.0005	0.34	ND	0.0093	0.12		
723-46-6	Sulfamethoxazole [µg/L]	0.000081	ND	ND	0.000028	ND	0.00026	0.00026	0.00017		
14808-79-8	Sulfate as SO4 [µg/L]	420000	100000	110000	220000	280000	140000	320000	240000		
18496-25-8	Sulfide [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
115-96-8	TCEP [µg/L]	0.0005	0.000026	ND	0.00018	ND	0.00054	ND	0.00019		
13674-84-5	TCPP [µg/L]	0.0026	0.0003	0.0003	0.0011	0.041	0.0018	ND	0.015		
13674-87-8	TDCPP [μg/L]	0.0017	ND	0.00014	0.00063	ND	0.0015	0.0034	0.0015		
TIC	Tentatively Identified Compounds [µg/L]	1.8	0.51	NA	0.82	ND	ND	9.7	2.9		
5902-51-2	Terbacil [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
13071-79-9	Terbufos [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
98-06-6	tert-Butylbenzene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
994-05-8	Tertiary-amyl methyl ether [µg/L]	ND	ND	ND	ND	ND	ND	ND	NA		
75-65-0	Tertiary-butyl alcohol [µg/L]	ND	0.35	ND	0.12	ND	ND	ND	ND		
58-22-0	Testosterone [µg/L]	0.000075	0.000015	ND	0.000032	ND	ND	ND	ND		
127-18-4	Tetrachloroethene (PCE) [µg/L]	27.7	0.77	1.5	10.4	54	0.39	41	31		
109-99-9	Tetrahydrofuran [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
479-45-8	Tetryl [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
7440-28-0	Thallium [μg/L]	0.0057	0.00087	0.00088	0.0026	ND	ND	ND	ND		
28249-77-6	Thiobencarb [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
THM	THMs, Total [μg/L]	1.2	0.55	0.59	0.81	3.7	0.66	0.9	1.8		
7440-29-1	Thorium [µg/L]	0.019	0.0048	0.0086	0.011	ND	0.064	ND	0.022		
ODOR	Threshold Odor Number [Units]	250	8.7	7.5	92	4	100	2	37		
108-88-3	Toluene [µg/L]	0.58	0.015	ND	0.21	ND	ND	ND	ND		
TDS	Total Dissolved Solids [µg/L]	790000	510000	500000	610000	990000	650000	1200000	930000		
TOC	Total Organic Carbon [μg/L]	550	910	39	520	1300	1400	510	1100		
1336-36-3	Total PCBs [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
TSuS	Total Suspended Solids [µg/L]	4100	130	10000	4600	54000	2000	200	20000		
1330-20-7	Total Xylenes [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		







			Using UCL9	95 for Monitoring	Well Data*	Using Maximum Value for 2020 Production Well Data**					
CASRN	Constituent [Units]	Estimated Con	ncentrations for Pr	oduction Wells	Plant Influent Concentration	Observed Con	centrations For P	roduction Well	Plant Influent Concentration		
		SM-4	SM-8	SM-9	Estimates	SM-4	SM-8	SM-9	Estimates		
8001-35-2	Toxaphene [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
TPHg	TPHg [μg/L]	ND	ND	NA	ND	ND	ND	ND	NA		
156-60-5	trans-1,2-Dichloroethene [µg/L]	ND	0.042	0.037	0.026	ND	ND	ND	ND		
10061-02-6	trans-1,3-Dichloropropene [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
110-57-6	trans-1,4-Dichloro-2-butene [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
24017-47-8	Triazofos [μg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
75-96-7	Tribromoacetic acid [μg/L]	ND	ND	ND	ND	ND	ND	ND	NA		
76-03-9	Trichloroacetic acid [µg/L]	ND	0.038	ND	0.013	ND	ND	ND	ND		
79-01-6	Trichloroethene (TCE) [µg/L]	22.2	0.92	0.44	8.2	59	7	0.34	23		
75-69-4	Trichlorofluoromethane [µg/L]	0.22	0.027	ND	0.088	ND	ND	ND	ND		
3380-34-5	Triclosan [μg/L]	ND	ND	ND	ND	0.0093	ND	ND	0.0033		
1582-09-8	Trifluralin [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
738-70-5	Trimethoprim [µg/L]	0.000087	0.000021	ND	0.000038	0.00054	0.00028	ND	0.00029		
786-19-6	Trithion [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
10028-17-8	Tritium [pci/L]	51	58	58	56	230	38	58	110		
TURB	Turbidity [ntu]	6.9	1.4	3.7	4	48	1.3	0.05	17		
7440-61-1	Uranium Rad [pci/L]	5.3	1.3	2.1	2.9	7.1	5.1	14	8.5		
7440-62-2	Vanadium, Total [μg/L]	2.4	0.13	0.58	1	4.2	6.1	3.5	4.7		
108-05-4	Vinyl Acetate [µg/L]	ND	ND	ND	ND	ND	ND	ND	ND		
75-01-4	Vinyl Chloride [μg/L]	ND	0.081	ND	0.028	ND	ND	ND	ND		
7440-66-6	Zinc [µg/L]	5.1	2.6	2.6	3.5	9.1	54	84	47		

## Notes:

CASRN = Chemical Abstracts Service Registry Number (as applicable);  $\mu g/L = micrograms$  per liter; mr = nanograms per liter; mr = nanogra units; 'ND' = non-detect; 'NA' = not available; '-' = not applicable.

\*UCL95 from statistical analysis of the monitoring well groundwater quality data set. In the event the UCL95 statistics could not be calculated, the maximum value was adopted as a conservative approach. For non-detect (ND) results a value of 0 was assumed \*\*Maximum observed values from production wells from available 2020 sampling data (provided for information only).







## Appendix G Mann-Kendall and Sen's Slope Trend Analysis Results







			Ana	lyses pe	rformed	on sample d	lata fr	om January 20	12 to June 2020 (inclusive	·)					
				Basic An	alysis			Mann-Kendall Trend Analysis							
Parameter	Count	Min (units)	Median (units)	Max (units)	Average (units)	Standard Deviation (units)	Units	P Value of Two Tailed Test (-)	Inferred Confidence Level (of Trend Present in Data Set) (%)	Slope (units/yr)	Normalized Slope (%/yr)	Meets Trend Assumptions?	Statistically Significant Mann-Kendall Trend		
CNA 2		(units)	(units)	(uiiits)	(units)	(units)	(-)	(-)	(70)	(units/yr)	(70/ yl)		Walli-Kelluali Heli		
SM-3	07	. 0.5	0.25	0.0	0.20	0.12	/1					N (lea - DDI data)			
1,1-Dichloroethene (1,1-DCE)	87	< 0.5	0.25	0.8 < 0.005	0.30 0.0025	0.13	μg/L					N (Ins. >RDL data)			
1,2,3-Trichloropropane (1,2,3-TCP)  1,4-Dioxane	7 29		1.40	7.9	2.22	0.0 2.13	μg/L	0.00	100.0%	0.20	21.42	N (Ins. trend data)	 De		
Aluminum	3	< 1.0 < 50	25.00	< 50	25.00	0.00	μg/L			-0.30	-21.42	N (Ins. trend data)	Down		
	3		0.70	< 2	0.73	0.00	μg/L								
Arsenic Barium		0.5 49.2	1	59.5	53.43	5.39	μg/L					N (Ins. trend data)			
	3		51.60				μg/L					N (Ins. trend data)			
Boron	0					0.00	μg/L					N (Ins. trend data)			
Carbon Tetrachloride (CTC)	87	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)			
Chromium	4	2.20	4.05	< 10	3.83	1.41	μg/L		<del></del>			N (Ins. trend data)			
cis-1,2-Dichloroethene (cis-1,2-DCE)	87	< 0.5	0.25	0.6	0.26	0.064	μg/L					N (Ins. > RDL data)			
Fluoride	3	330.00	370.00	390.00	363.33	30.55	μg/L					N (Ins. trend data)			
Gross Alpha	2	3.80	7.40	11.00	7.40	5.09	pci/L					N (Ins. trend data)			
Lead	3	< 5	2.50	< 5	2.50	0.00	μg/L					N (Ins. trend data)			
Manganese	3	4.60	5.50	6.70	5.60	1.05	μg/L					N (Ins. trend data)			
Nitrate (as N)	32	3800	4951	7341	5220	889	μg/L	0.07	93.1%	-136	-2.75	Υ			
Nitrate + Nitrite (as N)	2	440	2370	4300	2370	2729	μg/L					N (Ins. trend data)			
n-Nitrosodiethylamine (NDEA)	0						μg/L					N (Ins. trend data)			
Perchlorate	3	< 4	2.0	< 4	2.0	0.0	μg/L					N (Ins. trend data)			
Perfluorooctanoic Acid (PFOA)	1	0.0017	0.0017	0.0017	0.0017	0.00	μg/L					N (Ins. trend data)			
Perfluorooctanesulfonic Acid (PFOS)	1	0.0017	0.0017	0.0017	0.0017	0.00	μg/L					N (Ins. trend data)			
Selenium	3	2.7	2.7	< 5	2.8	0.3	μg/L					N (Ins. trend data)			
Tetrachloroethene (PCE)	87	5.2	8.40	26.6	10.43	4.12	μg/L	0.00	100.0%	-1.20	-14.32	Υ	Down		
Total Trihalomethanes (TTHMs)	12	0.55	1.05	2.8	1.25	0.69	μg/L	1.00	0.0%	0.00	0.00	Υ			
Trichloroethene (TCE)	87	< 0.5	3.30	8.7	3.88	1.97	μg/L	0.01	98.6%	-0.24	-7.36	Υ			
Uranium Rad	2	8	8.20	8.4	8.20	0.28	pci/L					N (Ins. trend data)			
Vanadium	2	4.4	4.70	5	4.70	0.42	μg/L					N (Ins. trend data)			
1,1,2-Trichloroethane (1,1,2-TCA)	12	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)			
1,1-Dichloroethane (1,1-DCA)	87	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)			
1,2-Dichloroethane (1,2-DCA)	87	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)			
Antimony	3	< 6	3.00	< 6	3.00	0.00	μg/L					N (Ins. trend data)			
Benzene	87	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)			
Bis(2-ethylhexyl)phthalate	6	< 3	1.50	< 3	1.50	0.00	μg/L					N (Ins. trend data)			
Bromate	0						μg/L					N (Ins. trend data)			
Hexachlorobenzene (HCB)	4	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. trend data)			
Methyl Tert-butyl Ether (MTBE)	87	< 3	1.50	< 3	1.50	0.00	μg/L					N (Ins. >RDL data)			
Methylene Chloride (also Dichloromethane [DCM])	87	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. > RDL data)			
Nickel	3	< 10	5.00	< 10	5.00	0.00	μg/L					N (Ins. trend data)			
Nitrite (as N)	3	< 400	200	< 400	200	0.00	μg/L					N (Ins. trend data)			
Strontium-90	0						μg/L pci/L					N (Ins. trend data)			
trans-1,2-Dichloroethene (trans-1,2-DCE)	87	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)			
trans-   /-  ) chioroethene (trans-   /-  )( -)															







			Anal	lyses pe	rformed o	on sample o	data fro	m January 20	112 to June 2020 (inclusive	·)						
				Basic An	alysis			Mann-Kendall Trend Analysis								
Parameter	Count	Min	Median	Max	Average	Standard Deviation	Units	P Value of Two Tailed Test	Inferred Confidence Level (of Trend Present in Data Set)	Slope	Normalized Slope	Meets Trend Assumptions?	Statistically Significant			
		(units)	(units)	(units)	(units)	(units)	(-)	(-)	(%)	(units/yr)	(%/yr)	Assumptions:	Mann-Kendall Trend			
SM-4	,															
1,1-Dichloroethene (1,1-DCE)	106	1.1	1.80	2.9	1.86	0.39	μg/L	0.00	100.0%	0.11	6.05	Υ				
1,2,3-Trichloropropane (1,2,3-TCP)	26	0.0026	0.0054	0.0100	0.0049	0.0023	μg/L					N (Ins. >RDL data)				
1,4-Dioxane	32	11.0	19.0	26.0	18.3	2.94	μg/L	0.00	100.0%	0.87	4.58	Υ				
Aluminum	4	3.4	16.2	< 50	15.2	11.4	μg/L					N (Ins. trend data)				
Arsenic	4	0.50	0.65	0.71	0.63	0.10	μg/L					N (Ins. trend data)				
Barium	4	56.5	61.8	76.0	64.0	8.5	μg/L					N (Ins. trend data)				
Boron	3	130	130	200	153	40.4	μg/L					N (Ins. trend data)				
Carbon Tetrachloride (CTC)	106	< 0.5	0.3	< 0.5	0.3	0.00	μg/L					N (Ins. >RDL data)				
Chromium	4	7.2	5.0	< 10	5.6	1.10	μg/L					N (Ins. trend data)				
cis-1,2-Dichloroethene (cis-1,2-DCE)	106	0.45	2.20	3.4	2.28	0.44	μg/L	0.00	100.0%	-0.10	-4.72	Υ				
Fluoride	4	290	335	350	328	26.3	μg/L					N (Ins. trend data)				
Gross Alpha	3	< 3	3.50	3.86	2.95	1.27	pci/L					N (Ins. trend data)				
Lead	4	4	2.50	< 5	2.88	0.75	μg/L					N (Ins. trend data)				
Manganese	5	4.17	25.10	33	22.67	10.90	μg/L					N (Ins. trend data)				
Nitrate (as N)	38	4653	6133	7400.0	6179	511	μg/L	0.90	10.0%	0.00	0.00	Υ				
Nitrate + Nitrite (as N)	3	570	5700	5800	4023	2991	μg/L					N (Ins. trend data)				
n-Nitrosodiethylamine (NDEA)	1	0.0015	0.0015	0.0015	0.0015	0.00	μg/L					N (Ins. trend data)				
Perchlorate	4	1.60	2.00	< 4	1.90	0.20	μg/L					N (Ins. trend data)				
Perfluorooctanoic Acid (PFOA)	2	0.0010	0.0013	0.0017	0.0013	0.0005	μg/L					N (Ins. trend data)				
Perfluorooctanesulfonic Acid (PFOS)	2	0.0017	0.0014	< 0.002	0.0014	0.0005	μg/L					N (Ins. trend data)				
Selenium	4	2.8	3.0	< 5	3.3	0.93	μg/L					N (Ins. trend data)				
Tetrachloroethene (PCE)	106	12.0	16.1	43.4	21.8	9.81	μg/L	0.00	100.0%	3.24	20.10	Υ	Up			
Total Trihalomethanes (TTHMs)	19	3.6	4.10	4.8	4.14	0.35	μg/L	0.65	35.4%	-0.025	-0.61	Υ				
Trichloroethene (TCE)	106	32.6	45.2	63.2	46.2	6.90	μg/L	0.98	1.5%	0.00	0.00	Υ				
Uranium Rad	3	5.9	6.0	7.1	6.3	0.67	pci/L					N (Ins. trend data)				
Vanadium	3	4.0	4.2	4.5	4.2	0.25	μg/L					N (Ins. trend data)				
1,1,2-Trichloroethane (1,1,2-TCA)	20	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)				
1,1-Dichloroethane (1,1-DCA)	106	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)				
1,2-Dichloroethane (1,2-DCA)	106	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)				
Antimony	4	0.2	3.0	< 6	2.3	1.40	μg/L					N (Ins. trend data)				
Benzene	106	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)				
Bis(2-ethylhexyl)phthalate	7	< 3	1.5	< 5	1.6	0.38	μg/L					N (Ins. trend data)				
Bromate	1	< 25	12.5	< 25	12.5	0.00	μg/L					N (Ins. trend data)				
Hexachlorobenzene (HCB)	5	< 0.5	0.25	< 1	0.30	0.11	μg/L					N (Ins. trend data)				
Methyl Tert-butyl Ether (MTBE)	105	< 2	1.5	< 3	1.5	0.05	μg/L					N (Ins. >RDL data)				
Methylene Chloride (also Dichloromethane [DCM])	106	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)				
Nickel	4	4.1	5.0	< 10	4.8	0.45	μg/L					N (Ins. trend data)				
Nitrite (as N)	4	< 100	200	< 400	163	75.00	μg/L					N (Ins. trend data)				
Strontium-90	1	< 0	0.0	< 0	0.0	0.00	pci/L					N (Ins. trend data)				
trans-1,2-Dichloroethene (trans-1,2-DCE)	106	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)				
Vinyl Chloride	19	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)				







Analyses performed on sample data from January 2017 to June 2020 (inclusive)													
				Basic An	alysis	-			Man	n-Kendall Tre	end Analysis		
Parameter	Count	Min	Median	Max	Average	Standard Deviation	Units	P Value of Two Tailed Test	Inferred Confidence Level (of Trend Present in Data Set)	Slope	Normalized Slope	Meets Trend Assumptions?	Statistically Significant
		(units)	(units)	(units)	(units)	(units)	(-)	(-)	(%)	(units/yr)	(%/yr)	7.00.000	Mann-Kendall Trend
SM-3													
1,1-Dichloroethene (1,1-DCE)	87	< 0.5	0.25	0.8	0.30	0.13	μg/L					N (Ins. >RDL data)	
1,2,3-Trichloropropane (1,2,3-TCP)	7	< 0.005	0.0025	< 0.005	0.0025	0.00	μg/L					N (Ins. trend data)	
1,4-Dioxane	29	< 1.0	1.40	7.9	2.22	2.13	μg/L	0.16	83.6%	0.17	15.53	Y	
Aluminum	1	< 50	25	< 50	25	0.00	μg/L					N (Ins. trend data)	
Arsenic	1	< 2	1.00	< 2	1.00	0.00	μg/L					N (Ins. trend data)	
Barium	1	59.5	59.5	59.5	59.5	0.00	μg/L					N (Ins. trend data)	
Boron	0						μg/L					N (Ins. trend data)	
Carbon Tetrachloride (CTC)	40	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
Chromium	1	< 10	5.00	< 10	5.00	0.00	μg/L					N (Ins. trend data)	
cis-1,2-Dichloroethene (cis-1,2-DCE)	87	< 0.5	0.25	0.6	0.26	0.064	μg/L					N (Ins. >RDL data)	
Fluoride	1	390	390	390	390	0.00	μg/L · "					N (Ins. trend data)	
Gross Alpha	1	3.8	3.8	3.8	3.8	0.00	pci/L					N (Ins. trend data)	
Lead	1	< 5	2.50	< 5	2.50	0.00	μg/L					N (Ins. trend data)	
Manganese	1	6.7	6.7	6.7	6.7	0.00	μg/L		24.10/		2.41	N (Ins. trend data)	
Nitrate (as N)	32	3800.0	4951	7341	5220	889	μg/L	0.66	34.1%	113	2.41	Y N. (Landan and Alada)	
Nitrate + Nitrite (as N)	1	4300	4300	4300	4300	0.00	μg/L					N (Ins. trend data)	
n-Nitrosodiethylamine (NDEA)	1		2.0		2.0	0.00	μg/L					N (Ins. trend data)	
Perchlorate Perfluorooctanoic Acid (PFOA)	1	< 4 0.0017	2.0 0.0017	< 4	2.0 0.0017	0.00	μg/L					N (Ins. trend data)	
Perfluorooctanoic Acid (PFOA)  Perfluorooctanesulfonic Acid (PFOS)	1	0.0017	0.0017	0.0017	0.0017	0.00	μg/L					N (Ins. trend data) N (Ins. trend data)	
Selenium	1	2.7	2.7	2.7	2.7	0.00	μg/L					N (Ins. trend data)	
Tetrachloroethene (PCE)	87	5.2	8.4	26.6	10.4	4.12	μg/L μg/L	0.16	84.2%	0.25	3.51	v	
Total Trihalomethanes (TTHMs)	12	0.55	1.05	2.8	1.25	0.69	μg/L μg/L	1.00	0.0%	0.23	20.65	N (Ins. trend data)	
Trichloroethene (TCE)	87	< 0.5	3.30	8.7	3.88	1.97	μg/L μg/L	0.01	99.3%	0.21	10.51	v	
Uranium Rad	1	8.40	8.40	8.40	8.40	0.00	μg/L pci/L			0.55		N (Ins. trend data)	
Vanadium	0						μg/L					N (Ins. trend data)	
1,1,2-Trichloroethane (1,1,2-TCA)	3	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. trend data)	
1,1-Dichloroethane (1,1-DCA)	40	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
1,2-Dichloroethane (1,2-DCA)	40	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
Antimony	1	< 6	3.00	< 6	3.00	0.00	μg/L					N (Ins. trend data)	
Benzene	40	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
Bis(2-ethylhexyl)phthalate	2	< 3	1.50	< 3	1.50	0.00	μg/L					N (Ins. trend data)	
Bromate	0						μg/L					N (Ins. trend data)	
Hexachlorobenzene (HCB)	1	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. trend data)	
Methyl Tert-butyl Ether (MTBE)	40	< 3	1.50	< 3	1.50	0.00	μg/L					N (Ins. >RDL data)	
Methylene Chloride (also Dichloromethane [DCM])	40	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
Nickel	1	< 10	5.00	< 10	5.00	0.00	μg/L					N (Ins. trend data)	
Nitrite (as N)	1	< 400	200	< 400	200	0.00	μg/L					N (Ins. trend data)	
Strontium-90	0						pci/L					N (Ins. trend data)	
trans-1,2-Dichloroethene (trans-1,2-DCE)	40	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
Vinyl Chloride	3	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. trend data)	







				,						,			
			Ana	lyses pe	rformed	on sample o	data fro	om January 20	17 to June 2020 (inclusive	?)			
				Basic An	alysis				Mar	nn-Kendall Tre	nd Analysis		
Parameter	Count	Min	Median	Max	Average	Standard Deviation	Units	P Value of Two Tailed Test	Inferred Confidence Level (of Trend Present in Data Set)	Slope	Normalized Slope	Meets Trend Assumptions?	Statistically Significant
		(units)	(units)	(units)	(units)	(units)	(-)	(-)	(%)	(units/yr)	(%/yr)		Mann-Kendall Trend
SM-4	106	4.4	4.00	2.0	4.06	0.20		0.00	100.00/	0.00	40.00		
1,1-Dichloroethene (1,1-DCE)	106	1.1	1.80	2.9	1.86	0.39	μg/L	0.00	100.0%	0.22	10.22	Y	
1,2,3-Trichloropropane (1,2,3-TCP)	26	0.0026	0.0054	0.0100	0.0049	0.00	μg/L					N (Ins. >RDL data)	
1,4-Dioxane	32	11.0	19.0	26.0	18.3	2.94	μg/L	0.01	99.4%	1.34	6.70	N (leaster and date)	Up
Aluminum	2	3.4	14.2	< 50	14.2	15.27	μg/L		<del></del>			N (Ins. trend data)	
Arsenic	2	0.6	0.7	0.71	0.7	0.08	μg/L		<del></del>			N (Ins. trend data)	
Barium	2	60.0	68.0	76.0	68.0	11.31	μg/L					N (Ins. trend data)	
Boron	2	130.0	130.0	130.0	130.0	0.00	μg/L					N (Ins. trend data)	
Carbon Tetrachloride (CTC)	43	< 0.5	0.3	< 0.5	0.3	0.00	μg/L					N (Ins. >RDL data)	
Chromium	2	7.2	6.1	< 10	6.1	1.56	μg/L					N (Ins. trend data)	
cis-1,2-Dichloroethene (cis-1,2-DCE)	106	0.45	2.20	3.4	2.28	0.44	μg/L	0.01	99.2%	-0.079	-3.95	Υ	
Fluoride	2	290.0	315.0	340.0	315.0	35.36	μg/L					N (Ins. trend data)	
Gross Alpha	2	< 3	2.50	3.5	2.50	1.41	pci/L					N (Ins. trend data)	
Lead	2	4	3.25	< 5	3.25	1.06	μg/L					N (Ins. trend data)	
Manganese	2	25.1	29.1	33.0	29.1	5.59	μg/L					N (Ins. trend data)	
Nitrate (as N)	38	4653	6133	7400.0	6179	511	μg/L	0.96	4.0%	0.00	0.00	Υ	
Nitrate + Nitrite (as N)	2	5700	5750	5800	5750	71	μg/L					N (Ins. trend data)	
n-Nitrosodiethylamine (NDEA)	1	0.0015	0.0015	0.0015	0.0015	0.00	μg/L					N (Ins. trend data)	
Perchlorate	2	1.6	2	< 4	1.8	0.28	μg/L					N (Ins. trend data)	
Perfluorooctanoic Acid (PFOA)	2	0.00096	0	0.0017	0.0	0.00	μg/L					N (Ins. trend data)	
Perfluorooctanesulfonic Acid (PFOS)	2	0.0017	0	< 0.002	0.0	0.00	μg/L					N (Ins. trend data)	
Selenium	2	3.1	4	4.6	3.9	1.06	μg/L					N (Ins. trend data)	
Tetrachloroethene (PCE)	106	12.0	16.1	43.4	21.8	9.81	μg/L	0.00	100.0%	6.39	18.62	Υ	Up
Total Trihalomethanes (TTHMs)	19	3.6	4.10	4.8	4.14	0.35	μg/L	0.73	26.6%	-0.20	-5.12	N (Ins. trend data)	
Trichloroethene (TCE)	106	32.6	45.2	63.2	46.2	6.90	μg/L	0.01	98.8%	2.02	4.34	Υ	
Uranium Rad	2	5.9	6.5	7.1	6.5	0.85	pci/L					N (Ins. trend data)	
Vanadium	1	4.2	4.2	4.2	4.2	0.00	μg/L					N (Ins. trend data)	
1,1,2-Trichloroethane (1,1,2-TCA)	5	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. trend data)	
1,1-Dichloroethane (1,1-DCA)	43	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
1,2-Dichloroethane (1,2-DCA)	43	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
Antimony	2	0.2	1.6	< 6	1.6	1.98	μg/L					N (Ins. trend data)	
Benzene	43	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
Bis(2-ethylhexyl)phthalate	3	< 3	1.5	< 5	1.8	0.58	μg/L					N (Ins. trend data)	
Bromate	1	< 25	12.5	< 25	12.5	0.00	μg/L					N (Ins. trend data)	
Hexachlorobenzene (HCB)	2	< 0.5	0.38	< 1	0.38	0.18	μg/L					N (Ins. trend data)	
Methyl Tert-butyl Ether (MTBE)	43	< 2	1.5	< 3	1.5	0.08	μg/L					N (Ins. >RDL data)	
Methylene Chloride (also Dichloromethane [DCM])	43	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
Nickel	2	4.1	4.6	< 10	4.6	0.64	μg/L					N (Ins. trend data)	
Nitrite (as N)	2	< 100	125.0	< 400	125.0	106.07	μg/L					N (Ins. trend data)	
Strontium-90	1	< 0	0.0	< 0	0.0	0.00	pci/L					N (Ins. trend data)	
trans-1,2-Dichloroethene (trans-1,2-DCE)	43	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. >RDL data)	
Vinyl Chloride	4	< 0.5	0.25	< 0.5	0.25	0.00	μg/L					N (Ins. trend data)	







### Notes:

- 1. Non-detect multiplier of 0.5 applied to <DL sample data
- 2. **Bold** text in the 'Parameter' column indicates one or more analyses returned notable results
- 3. Basic analysis performed with the following conditions:
- Calculation of median and average used multiplier of 0.5 for <DL sample data
- 4. Mann-Kendall trend analysis:
- Bold text/values in the 'Mann-Kendall Trend Analysis' columns indicates a Mann-Kendall trend has been detected
- Statistically significant trend defined as:
- -- P-value of two-tailed test less than or equal to 0.10
- -- Absolute value of Sen's normalized slope is greater than 5.0%
- -- Absolute slope criteria met for individual parameters:
- 1,1-Dichloroethene: 0.5 μg/L/yr
- 1,4-Dioxane: 0.1 μg/L/yr
- cis-1,2-Dichloroethene: 0.5 μg/L/yr
- Nitrate as N: 200.0 μg/L/yr
- Tetrachloroethene: 0.5 μg/L/yr
- Total Trihalomethanes: 5.0 μg/L/yr
- Trichloroethene: 0.5 μg/L/yr
- Up: Denotes statistically significant upward trend
- Down: Denotes statistically significant downward trend
- "---": Denotes no trend
- Non-detect multiplier of 0.5 applied to <DL sample data
- 5. Legend for trend assumptions:
- Y: All conditions are met
- N: Not all conditions are met, specified as follows:
- -- No new data: No sample data present for years reported
- -- Ins. >DL data: Less than 50% of data was above detection limit
- -- Non-Monotonic: Trend is non-monotonic (i.e., trend reversal or termination is present)
- -- Damaged: The well is damaged
- -- Ins. trend data: Found fewer than 8 data points in the data set used to calculate Mann-Kendall results

Full Raw Water Quality Characterization - Olympic Well Field - Step 2 of 97-005 Evaluation Current Rev No.: 1: 308038-03533-00-EN-REP-0004







## Appendix H Trend Charts

#### - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) — Monthly Total Pumped (Million Gallons [MG]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) 1,1-Dichloroethene (1,1-DCE) 1,2,3-Trichloropropane (1,2,3-TCP) 1,4-Dioxane 10 42 10 0.1 0.09 35**(9M)** padmnd 21 d 35 28 21 14 7 7 Wonthly Total Pumped (MG) 0.08 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.07 0.06 0.05 Total 1 0.04 3 0.03 0.02 0.01 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 **Aluminum** Arsenic **Barium** 2500 1000 10 2250 900 35 28 21 14 7 7 Wonthly Total Pumped (MG) 28 28 21 Lotal Brumbed (MG) 2000 800 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 1750 700 1500 600 1250 500 1000 400 750 300 500 200 250 100 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 **Boron** Carbon Tetrachloride (CTC) Chromium 1000 42 50 42 900 45 0.9 35 (SW) padwnd 21 d 35 28 21 14 14 T 7 35 28 21 14 7 7 Wonthly Total Pumped (MG) 800 8.0 40 Concentration (µg/L) Concentration (µg/L) No Boron data for SM-03 Concentration (µg/L) 700 0.7 35 600 0.6 30 500 0.5 25 400 20 0.4 15 300 0.3 200 0.2 10 100 0.1 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020

**Appendix H - Production Well SM-03** 

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#### - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) — Monthly Total Pumped (Million Gallons [MG]) Maximum Contaminant Level (MCL) - Triangle symbols ((\( \text{\( \text{\( \)}\)}\)) indicate that values were converted from Nitrate (as NO3) to Nitrate (as N) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) cis-1,2-Dichloroethene (cis-1,2-DCE) **Fluoride Gross Alpha** 10 42 2000 1800 35**(M)** padmnd 21**d** 35 **(9W)** padwnd 1 35 M2 28 21 **Lotal Pumped (MG)** 1600 Concentration (pCi/L) Concentration (µg/L) Concentration (µg/L) 1400 1200 1000 14 P Total 14P 800 600 Monthly 400 200 2016 2017 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2018 2019 2020 Lead Manganese Nitrate (as N) 500 10000 42 42 14 13 12 450 9000 35 28 21 14 7 7 Wonthly Total Pumped (MG) 400 8000 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 28 21 21 **Lotal Pumped (** 350 7000 10 300 6000 250 5000 200 4000 150 3000 100 2000 50 1000 2012 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2012 2013 2014 2015 2016 2017 2018 2019 2020 2019 2020 Nitrate + Nitrite (as N) n-Nitrosodiethylamine (NDEA) Perchlorate 10000 0.1 42 10 42 9000 0.09 35 (M) padund 1 9 35 28 21 14 7 7 7 14 Total Pumbed (MG) 35 **S** 8000 0.08 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 28 21 14 7 7 Wonthly Total Pumped 7000 0.07 No n-Nitrosodiethylamine data for 6000 0.06 SM-03 5000 0.05 Monthly Total F 4000 0.04 3000 0.03 2000 0.02 1000 2012 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017

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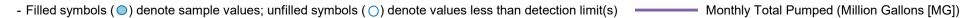
#### - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) Monthly Total Pumped (Million Gallons [MG]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) Perfluorooctanoic Acid (PFOA) Perfluorooctanesulfonic Acid (PFOS) Selenium 50 42 0.1 42 42 0.1 45 0.09 0.09 35 28 21 14 7 Wonthly Total Pumped (MG) 28 21 21 Lotal Bumped (MG) Monthly Total Pumped (MG) 35 0.08 0.08 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.07 0.07 28 30 0.06 0.06 25 21 0.05 0.05 0.04 0.04 20 15 0.03 0.03 7 Monthly 0.02 0.02 0.01 0.01 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 **Tetrachloroethene (PCE) Total Trihalomethanes (TTHMs)** Trichloroethene (TCE) 42 42 42 50 80 10 45 70 35 M2 28 21 Lotal Pumbed (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) Monthly Total Pumped (MG) 35 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 60 28 50 30 25 40 20 30 20 Monthly 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2012 2013 2014 2015 2018 2019 2020 **Uranium Rad** Vanadium 1,1,2-Trichloroethane (1,1,2-TCA) 20 42 50 42 42 18 45 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) Monthly Total Pumped (MG) 35 Concentration (pCi/L) Concentration (µg/L) Concentration (µg/L) 35 28 12 30 10 25 21 20 15 10 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020

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## - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) Monthly Total Pumped (Million Gallons [MG]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) 1,1-Dichloroethane (1,1-DCA) 1,2-Dichloroethane (1,2-DCA) Antimony 42 10 5 0.9 35 **(9W)** padwnd I 35**(9)** padwnd 1 35 28 21 14 7 7 Wonthly Total Pumped (MG) 8.0 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.7 0.6 0.5 14 7 7 Monthly Total F Total 14L 0.3 0.1 2017 2018 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2019 2020 2012 2013 Bis(2-ethylhexyl)phthalate (DEHP) Benzene **Bromate** 42 10 0.9 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) No Bromate data for SM-03 0.3 0.1 2012 2013 2014 2015 2016 2017 2018 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2019 2020 Hexachlorobenzene (HCB) Methyl Tert-butyl Ether (MTBE) Methylene Chloride (also Dichloromethane [DCM]) 15 14 42 42 0.9 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) 13 35 28 21 14 7 7 Wonthly Total Pumped (MG) Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.6 0.5 0.2 0.1 2013 2014 2015 2016 2017 2018 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020

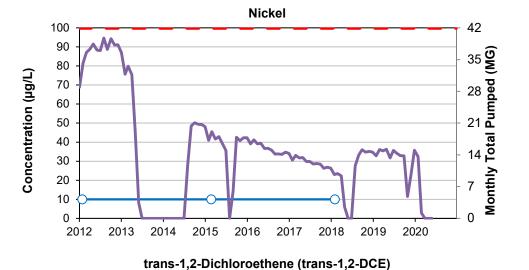
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# **Appendix H - Production Well SM-03**



42

35 28 21 14 7 7 Wonthly Total Pumped (MG)

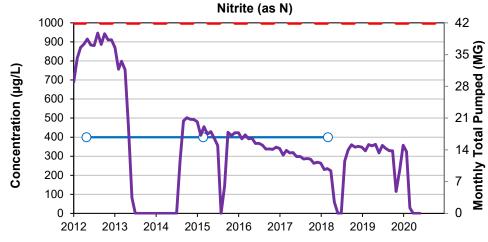


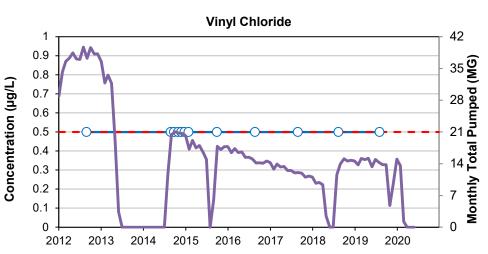
2012 2013 2014 2015 2016 2017 2018 2019 2020

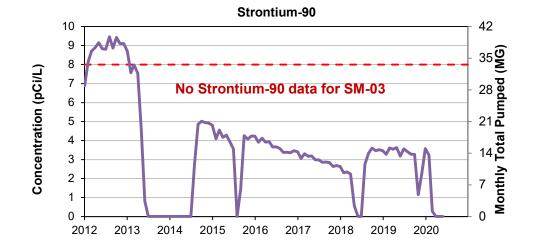
10

9

Concentration (µg/L)







Maximum Contaminant Level (MCL)

Secondary Maximum Contaminant Level (SMCL)

Notification Level (NL)

#### - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) Monthly Total Pumped (Million Gallons [MG]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) 1,1-Dichloroethene (1,1-DCE) 1,2,3-Trichloropropane (1,2,3-TCP) 1,4-Dioxane 42 10 0.02 42 30 27 0.018 35**(9M)** padmnd 21 d 28 21 21 Lotal Bumped (MG) 35 M2 28 21 Lotal Pumbed (MG) 0.016 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.014 21 0.012 0.01 Total 14P 12 0.008 0.006 3 onthly. Monthly . 0.004 0.002 2012 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 **Aluminum** Arsenic **Barium** 2500 1000 10 2250 900 35 **(9W)** pad md 21 d 35 28 21 14 7 7 Wonthly Total Pumped (MG) 2000 800 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 1750 700 1500 600 1250 500 1000 400 750 300 500 200 250 100 2012 2013 2014 2015 2016 2017 2018 2019 2020 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 **Boron** Carbon Tetrachloride (CTC) Chromium 1000 42 50 42 900 0.9 45 35 **(9W)** padund I 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) 8.0 800 40 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 35 700 0.7 600 30 0.6 500 0.5 25 **14 D** 20 400 0.4 15 300 0.3 200 10 0.2 100 0.1 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019

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#### - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) — Monthly Total Pumped (Million Gallons [MG]) Maximum Contaminant Level (MCL) - Triangle symbols (A) indicate that values were converted from Nitrate (as NO3) to Nitrate (as N) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) cis-1,2-Dichloroethene (cis-1,2-DCE) **Fluoride Gross Alpha** 10 42 2000 1800 35 **(SW)** padwn**d** 1 35**(9)** padwn**d** 21 **d** 1600 Concentration (pCi/L) Concentration (µg/L) Concentration (µg/L) 1400 1200 1000 14 L Total 800 600 Monthly 400 200 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 Lead Manganese Nitrate (as N) 500 10000 14 13 450 9000 35 (9M) padund I 28 21 21 **Lotal Pumped (MG)** 35 28 21 14 7 7 Wonthly Total Pumped (MG) 400 8000 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 350 7000 300 6000 250 5000 200 4000 150 3000 Monthly -100 2000 50 1000 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 Nitrate + Nitrite (as N) n-Nitrosodiethylamine (NDEA) Perchlorate 10000 0.1 42 10 42 9000 0.09 35 (SW) padwnd 1 21 d 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) 8000 0.08 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 7000 0.07 6000 0.06 5000 0.05 **1**4 **L** 4000 0.04 3000 0.03 2000 0.02 1000 2012 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019

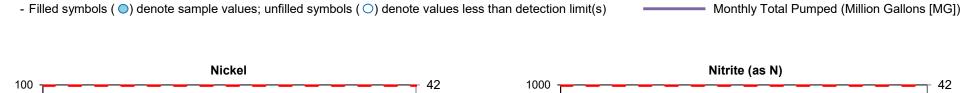
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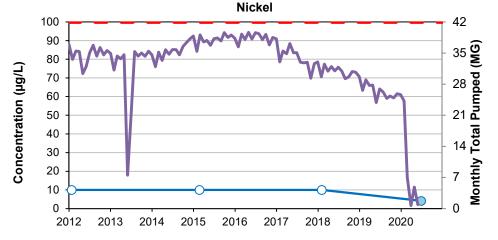
#### - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) Monthly Total Pumped (Million Gallons [MG]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) Perfluorooctanoic Acid (PFOA) Perfluorooctanesulfonic Acid (PFOS) Selenium 42 50 0.1 0.1 45 0.09 0.09 35**(9M)** padmnd 21 d 28 21 21 Lotal Bumped (MG) 35 M2 28 21 Lotal Pumbed (MG) 0.08 0.08 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.07 0.07 35 30 0.06 0.06 25 0.05 0.05 Total 1 20 0.04 0.04 0.03 0.03 15 Monthly . 0.02 0.02 0.01 0.01 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 **Tetrachloroethene (PCE) Total Trihalomethanes (TTHMs)** Trichloroethene (TCE) 60 80 70 55 70 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 21 Lotal Pumped (MG) 50 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 40 50 35 30 40 30 25 30 20 20 15 20 10 10 10 2018 2019 2020 2017 2012 2013 2014 2015 2016 2017 2013 2014 2015 2016 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 **Uranium Rad** Vanadium 1,1,2-Trichloroethane (1,1,2-TCA) 20 50 42 42 18 45 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) 40 Concentration (pCi/L) Concentration (µg/L) Concentration (µg/L) 35 12 30 10 25 20 10 2013 2014 2015 2016 2017 2018 2019 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018

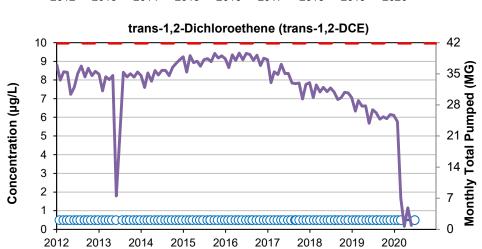
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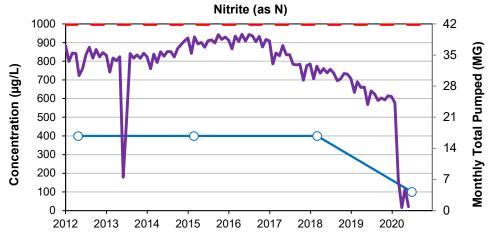
## - Filled symbols ( ○) denote sample values; unfilled symbols ( ○) denote values less than detection limit(s) —— Monthly Total Pumped (Million Gallons [MG]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) 1,1-Dichloroethane (1,1-DCA) 1,2-Dichloroethane (1,2-DCA) Antimony 42 0.9 35**(M)** padund 1 35 M2 28 21 Lotal Bumped (MG) 8.0 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.7 0.6 0.5 Total 1 0.4 0.3 Monthly . 0.1 2012 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 Benzene Bis(2-ethylhexyl)phthalate (DEHP) **Bromate** 30 42 30 27 27 0.9 35 28 21 4 Lotal Brumbed (MG) 35 (9M) padund I 35 28 21 14 7 7 Wonthly Total Pumped (MG) 8.0 24 24 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 21 18 18 15 12 12 0.3 Monthly -0.2 0.1 2018 2019 2020 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2014 2015 2016 2017 2013 2014 2015 2016 2017 2018 2013 Hexachlorobenzene (HCB) Methyl Tert-butyl Ether (MTBE) Methylene Chloride (also Dichloromethane [DCM]) 42 42 0.9 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) 35 28 21 14 7 7 Wonthly Total Pumped (MG) Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.6 0.5 0.4 0.2 0.1 2013 2014 2015 2016 2017 2018 2019 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018

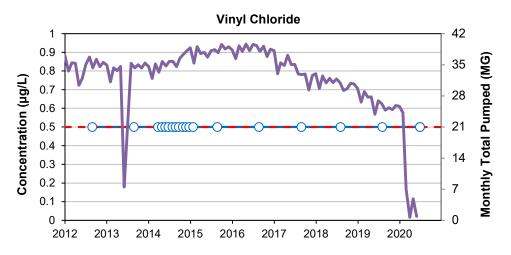
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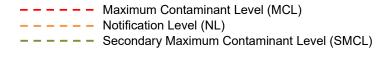


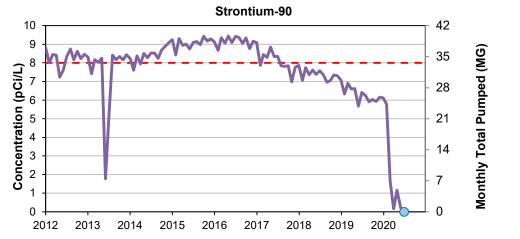


















# Appendix I COPC Concentration v. Precipitation Charts

#### Appendix I - Production Well SM-03 COPC Concentration v. Precipitation Charts - Filled symbols ( ○) denote sample values; unfilled symbols ( ○) denote values less than detection limit(s) Monthly Total Precipitation (inches [in]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) 1,1-Dichloroethene (1,1-DCE) 1,2,3-Trichloropropane (1,2,3-TCP) 1,4-Dioxane 10 10 Total Precipitation (in) 0.1 10 8 6 4 5 onthly Total Precipitation (in) 9 0.09 0.08 8 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.07 0.06 6 0.05 0.04 3 0.03 0.02 0.01 2014 2015 2016 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2012 2013 **Aluminum Arsenic Barium** 2500 1000 10 Monthly Total Precipitation (in) 2250 900 2000 800 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 1750 700 1500 6 600 1250 500 400 1000 750 300 500 200 250 100 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2013 2014 2015 2016 2017 2018 2019 2020 **Boron** Carbon Tetrachloride (CTC) Chromium 1000 Monthly Total Precipitation (in) 50 Monthly Total Precipitation (in) 900 45 0.9 40 800 8.0 Concentration (µg/L) No Boron data for SM-03 Concentration (µg/L) 0.7 35 700 600 0.6 30 500 25 400 20 0.4 300 15 0.3 10 200 0.2 100 0.1

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#### - Filled symbols ( ○) denote sample values; unfilled symbols ( ○) denote values less than detection limit(s) Monthly Total Precipitation (inches [in]) Maximum Contaminant Level (MCL) - Triangle symbols (A) indicate that values were converted from Nitrate (as NO3) to Nitrate (as N) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) cis-1,2-Dichloroethene (cis-1,2-DCE) **Fluoride Gross Alpha** 10 2000 Total Precipitation (in) Total Precipitation (in) 9 1800 13 1600 8 Concentration (pCi/L) Concentration (µg/L) 11 1400 1200 6 1000 800 3 600 lonthly . 400 200 2014 2015 2016 2017 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2013 2018 2019 Lead Manganese Nitrate (as N) 9 8 0 Precipitation (in) 500 10 8 6 4 2 0 Monthly Total Precipitation (in) 10000 14 13 450 9000 12 400 8000 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 350 7000 10 300 6 6000 250 5000 **Monthly Total** 200 4000 150 3000 100 2000 50 1000 2017 2018 2017 2018 2019 2017 2013 2014 2015 2016 2019 2020 2012 2013 2014 2015 2016 2012 2013 2014 2015 2016 2018 2019 2020 Nitrate + Nitrite (as N) n-Nitrosodiethylamine (NDEA) Perchlorate 10000 9 8 01 Precipitation (in) 0.1 9 8 01 Precipitation (in) 10 9000 0.09 8000 0.08 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 7000 0.07 No n-Nitrosodiethylamine data for 6000 0.06 SM-03 5000 0.05 4 2 0 Monthly Total F 4000 0.04 3000 0.03 Monthly. 2000 0.02 2 1000

Appendix I - Production Well SM-03 COPC Concentration v. Precipitation Charts

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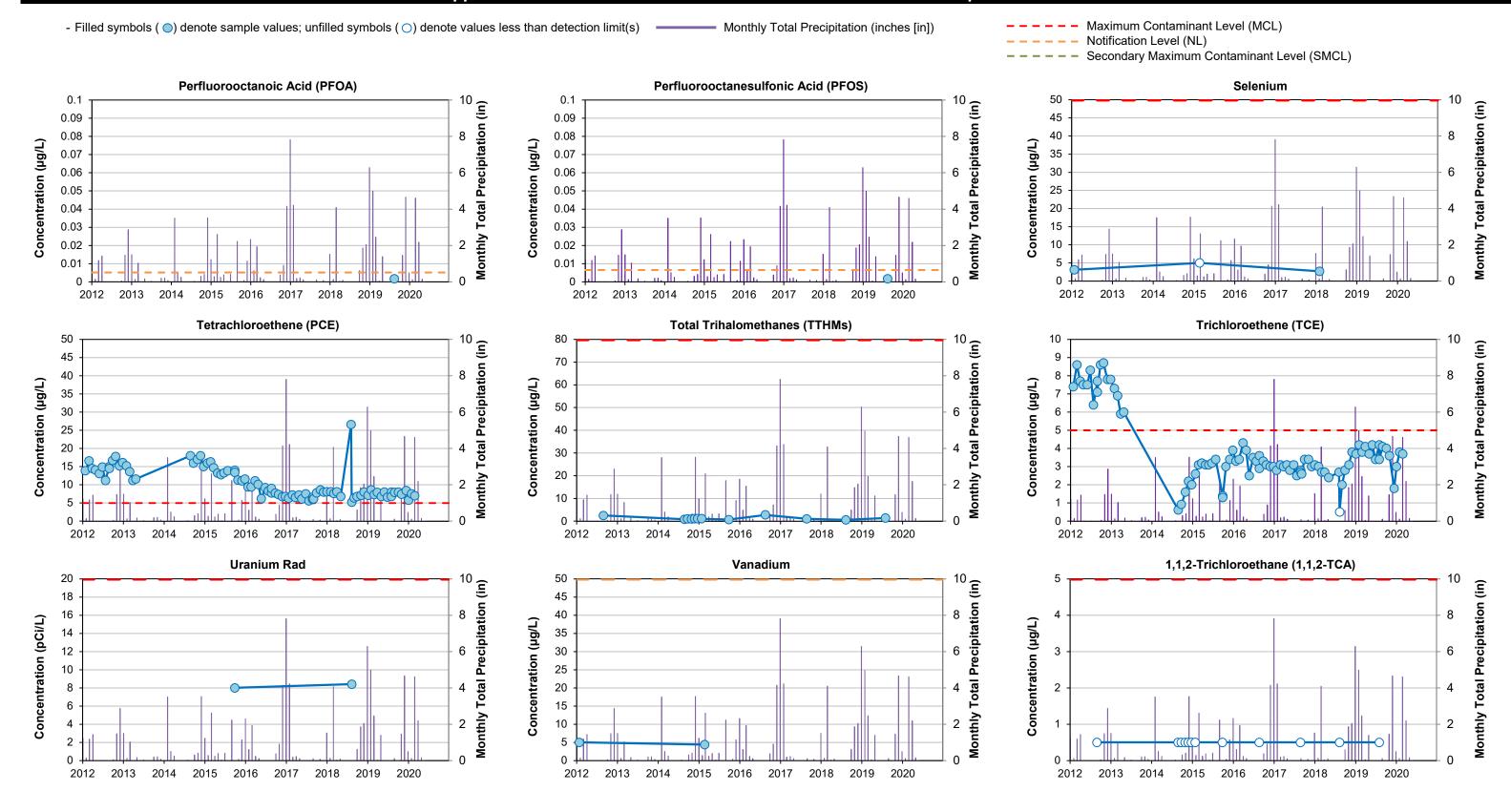
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## **Appendix I - Production Well SM-03 COPC Concentration v. Precipitation Charts**

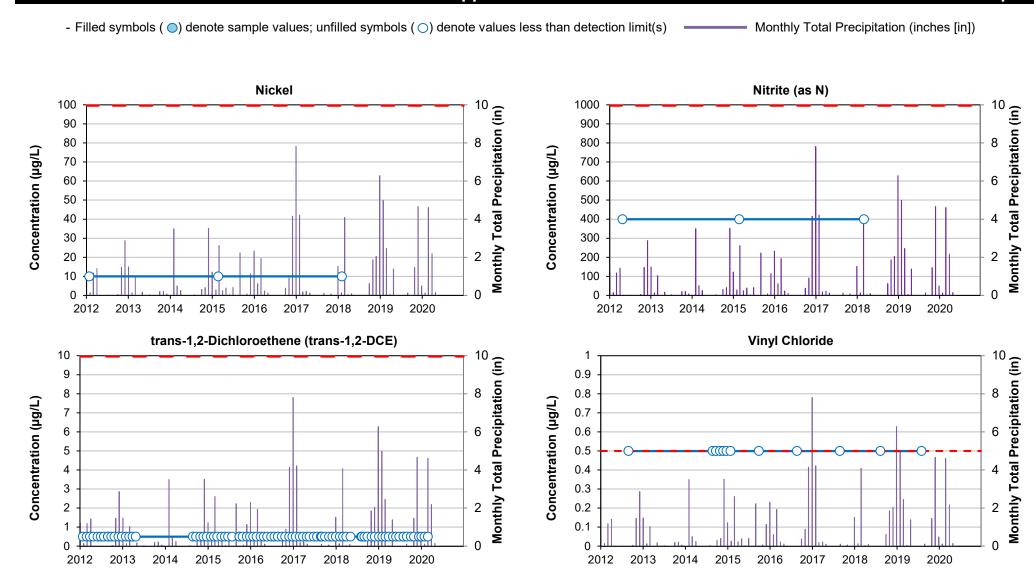


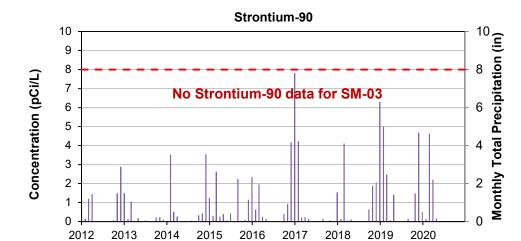
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## - Filled symbols ( ○) denote sample values; unfilled symbols ( ○) denote values less than detection limit(s) Monthly Total Precipitation (inches [in]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) 1,1-Dichloroethane (1,1-DCA) 1,2-Dichloroethane (1,2-DCA) Antimony 10 5 onthly Total Precipitation (in) 10 8 6 4 2 0 Monthly Total Precipitation (in) 0.9 8.0 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.7 0.6 6 0.3 0.1 2015 2013 2014 2017 2018 2019 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2016 2012 2013 Bis(2-ethylhexyl)phthalate (DEHP) Benzene **Bromate** 10 8 6 4 2 0 Monthly Total Precipitation (in) 10 8 6 4 2 0 Monthly Total Precipitation (in) 10 0.9 0.8 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) No Bromate data for SM-03 6 0.3 0.1 2015 2016 2017 2018 2014 2015 2016 2017 2018 2019 2020 2015 2016 2017 2018 2012 2013 2014 2019 2020 2013 2013 2014 Hexachlorobenzene (HCB) Methyl Tert-butyl Ether (MTBE) Methylene Chloride (also Dichloromethane [DCM]) 9 8 01 Precipitation (in) 10 8 6 4 2 0 Monthly Total Precipitation (in) 0.9 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.6 0.5 0.3 Monthly 0.2 2013 2014 2015 2016 2017 2018 2014 2015 2016 2017 2018 2013 2014 2015 2016 2017 2018

Appendix I - Production Well SM-03 COPC Concentration v. Precipitation Charts

# Appendix I - Production Well SM-03 COPC Concentration v. Precipitation Charts





– Maximum Contaminant Level (MCL)

Secondary Maximum Contaminant Level (SMCL)

Notification Level (NL)

#### Appendix I - Production Well SM-04 COPC Concentration v. Precipitation Charts - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) — Monthly Total Precipitation (inches [in]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) 1,1-Dichloroethene (1,1-DCE) 1,2,3-Trichloropropane (1,2,3-TCP) 1,4-Dioxane 10 0.02 30 Total Precipitation (in) 10 8 6 4 2 0 Monthly Total Precipitation (in) 27 9 0.018 0.016 24 Concentration (µg/L) Concentration (µg/L) 0.014 21 0.012 6 0.01 0.008 0.006 0.004 0.002 2015 2017 2019 2012 2013 2014 2015 2016 2017 2013 2014 2015 2016 2017 2018 2019 2020 2013 2014 2016 2018 2020 2018 2019 2020 **Aluminum** Arsenic **Barium** 2500 10 Monthly Total Precipitation (in) 1000 2250 900 2000 800 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 1750 700 1500 600 6 1250 500 1000 400 750 300 500 200 250 100 2012 2013 2014 2015 2016 2017 2018 2019 2020 2014 2015 2016 2017 2018 2012 2013 2014 2015 2016 2017 2018 2019 2020 2013 2019 **Boron** Carbon Tetrachloride (CTC) Chromium 1000 onthly Total Precipitation (in) Monthly Total Precipitation (in) 50 900 45 0.9 40 800 8.0 Concentration (µg/L) Concentration (µg/L) 0.7 35 700 600 30 0.6 500 0.5 25 20 400 300 15 0.3 200 10 0.2 100

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#### **Appendix I - Production Well SM-04 COPC Concentration v. Precipitation Charts** - Filled symbols ( ○) denote sample values; unfilled symbols ( ○) denote values less than detection limit(s) Monthly Total Precipitation (inches [in]) Maximum Contaminant Level (MCL) - Triangle symbols ( $\triangle$ ) indicate that values were converted from Nitrate (as NO3) to Nitrate (as N) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) cis-1,2-Dichloroethene (cis-1,2-DCE) **Fluoride Gross Alpha** 10 2000 9 1800 13 1600 Concentration (pCi/L) Concentration (µg/L) 1400 1200 6 1000 800 600 400 200 2017 2013 2014 2015 2016 2017 2018 2019 2020 2014 2016 2017 2018 2019 2020 2013 2014 2015 2016 2018 2019 2013 2015 Lead Manganese Nitrate (as N) 500 10 8 6 4 2 0 Monthly Total Precipitation (in) 10000 14 13 450 9000 12 400 8000 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 350 7000 10 300 6 6000 250 5000 **Monthly Total** 200 4000 150 3000 100 2000 50 1000 2018 2019 2020 2017 2018 2017 2012 2013 2014 2015 2016 2017 2013 2014 2015 2016 2019 2020 2012 2013 2014 2015 2016 2018 2019 2020 Nitrate + Nitrite (as N) n-Nitrosodiethylamine (NDEA) Perchlorate 10000 0.1 9 8 0 Precipitation (in) 10 9000 0.09 8000 0.08 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 7000 0.07 6000 0.06 5000 0.05 4 2 0 Monthly Total F 4000 0.04 3000 0.03 onthly 2000 0.02 1000

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#### **Appendix I - Production Well SM-04 COPC Concentration v. Precipitation Charts** - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) Monthly Total Precipitation (inches [in]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) Perfluorooctanoic Acid (PFOA) Perfluorooctanesulfonic Acid (PFOS) Selenium 0.1 Total Precipitation (in) 50 0.1 10 8 6 4 2 0 Monthly Total Precipitation (in) 45 0.09 0.09 0.08 0.08 40 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.07 0.07 35 30 0.06 0.06 6 25 0.05 0.05 20 0.04 0.04 0.03 0.03 15 0.02 0.02 0.01 0.01 2012 2013 2014 2015 2016 2017 2018 2017 2018 2019 2020 2012 2013 2014 2015 2016 2017 2018 2019 2020 2019 2020 2012 2013 2014 2015 2016 **Tetrachloroethene (PCE) Total Trihalomethanes (TTHMs)** Trichloroethene (TCE) 60 80 Monthly Total Precipitation (in) 70 55 70 50 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 45 60 40 50 35 30 40 30 25 30 20 20 15 20 10 2013 2014 2015 2017 2018 2019 2020 2015 2017 2018 2019 2020 2014 2015 2016 2017 2018 2019 2020 2016 2014 2016 2013 **Uranium Rad** Vanadium 1,1,2-Trichloroethane (1,1,2-TCA) 20 10 8 6 4 2 0 Monthly Total Precipitation (in) 50 Monthly Total Precipitation (in) 18 45 40 Concentration (pCi/L) Concentration (µg/L) Concentration (µg/L) 35 14 12 30 10 25 20 10

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### **Appendix I - Production Well SM-04 COPC Concentration v. Precipitation Charts** - Filled symbols ( ) denote sample values; unfilled symbols ( ) denote values less than detection limit(s) — Monthly Total Precipitation (inches [in]) Maximum Contaminant Level (MCL) Notification Level (NL) Secondary Maximum Contaminant Level (SMCL) 1,1-Dichloroethane (1,1-DCA) 1,2-Dichloroethane (1,2-DCA) Antimony onthly Total Precipitation (in) 10 8 6 4 2 0 Monthly Total Precipitation (in) 10 5 0.9 8.0 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.7 0.6 6 0.5 0.3 0.1 2013 2014 2015 2017 2018 2019 2014 2015 2016 2017 2018 2019 2020 2013 2014 2015 2016 2017 2018 2019 2020 2016 2012 2013 Bis(2-ethylhexyl)phthalate (DEHP) Benzene **Bromate** 30 10 8 6 4 2 0 Monthly Total Precipitation (in) 10 8 6 4 2 0 Monthly Total Precipitation (in) 30 27 0.9 27 0 0.8 24 24 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 21 18 6 18 15 0.3 0.1 2015 2016 2017 2018 2014 2015 2016 2017 2018 2019 2020 2015 2017 2018 2012 2013 2014 2019 2020 2013 2013 2014 2016 Hexachlorobenzene (HCB) Methyl Tert-butyl Ether (MTBE) Methylene Chloride (also Dichloromethane [DCM]) . 0 I Precipitation (in) Monthly Total Precipitation (in) 0.9 Concentration (µg/L) Concentration (µg/L) Concentration (µg/L) 0.6 0.5 0.3 Monthly . 0.2 0.1

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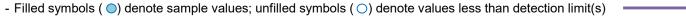
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## Appendix I - Production Well SM-04 COPC Concentration v. Precipitation Charts



Monthly Total Precipitation (inches [in])

