

**CLIENT NAME: DEFENCE CONSTRUCTION CANADA  
14 ALERT BLVD, 8 WING TRENTON  
ASTRA, ON K0K3W0  
(613) 392-2811**

**ATTENTION TO: Cameron Chadwick**

**PROJECT: Eureka Water License**

**AGAT WORK ORDER: 18T358833**

**MICROBIOLOGY ANALYSIS REVIEWED BY: Nivine Basily, Inorganics Report Writer**

**TRACE ORGANICS REVIEWED BY: Oksana Gushyla, Trace Organics Lab Supervisor**

**WATER ANALYSIS REVIEWED BY: Parvathi Malemath, Data Reviewer**

**DATE REPORTED: Jul 17, 2018**

**PAGES (INCLUDING COVER): 15**

**VERSION\*: 1**

Should you require any information regarding this analysis please contact your client services representative at (905) 712-5100

**\*NOTES**

**All samples will be disposed of within 30 days following analysis. Please contact the lab if you require additional sample storage time.**



# Certificate of Analysis

AGAT WORK ORDER: 18T358833

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5835 COOPERS AVENUE  
MISSISSAUGA, ONTARIO  
CANADA L4Z 1Y2  
TEL (905)712-5100  
FAX (905)712-5122  
<http://www.agatlabs.com>

CLIENT NAME: DEFENCE CONSTRUCTION CANADA

SAMPLING SITE:

ATTENTION TO: Cameron Chadwick

SAMPLED BY:

## Microbiological Analysis (water)

DATE RECEIVED: 2018-07-06

DATE REPORTED: 2018-07-17

Parameter	Unit	SAMPLE DESCRIPTION:		ERK-4	ERK-5	ERK-1
		SAMPLE TYPE:	DATE SAMPLED:	Water	Water	Water
		G / S	DATE SAMPLED:	2018-07-01	2018-07-01	2018-07-01
Fecal Coliform	CFU/100mL	2	RDL	9381766	9381776	RDL 9381778

Comments: RDL - Reported Detection Limit; G / S - Guideline / Standard

9381766-9381778 RDL >1 indicates dilutions of the sample.

ND - Not Detected.

HOLDING TIME-the time from sample collection to initiation of analysis exceeded 48 hours



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SAMPLED BY:

## Oil and Grease (+Total) - water

DATE RECEIVED: 2018-07-06

DATE REPORTED: 2018-07-17

Parameter	Unit	SAMPLE DESCRIPTION:		ERK-4	ERK-5	ERK-1
		G / S	RDL	SAMPLE TYPE:	Water	Water
				DATE SAMPLED:	2018-07-01	2018-07-01
Oil and Grease (animal/vegetable)	mg/L			0.5	0.8	1.2
Oil and Grease (mineral) in water	mg/L			0.5	<0.5	<0.5
Oil and Grease (Total) in water	mg/L			0.5	1.1	1.3
						1.4

Comments: RDL - Reported Detection Limit; G / S - Guideline / Standard

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## PHCs F1 - F4 (Water)

DATE RECEIVED: 2018-07-06

DATE REPORTED: 2018-07-17

Parameter	Unit	SAMPLE DESCRIPTION:		ERK-4	ERK-5
		SAMPLE TYPE:		Water	Water
		G / S	DATE SAMPLED:	2018-07-01	2018-07-01
Benzene	µg/L	370	0.20	<0.20	<0.20
Toluene	µg/L	2	0.20	<0.20	<0.20
Ethylbenzene	µg/L	90	0.10	<0.10	<0.10
Xylene Mixture	µg/L		0.20	<0.20	<0.20
F1 (C6 - C10)	µg/L		25	<25	<25
F1 (C6 to C10) minus BTEX	µg/L		25	<25	<25
F2 (C10 to C16)	µg/L		100	<100	<100
F3 (C16 to C34)	µg/L		100	<100	<100
F4 (C34 to C50)	µg/L		100	<100	<100
Gravimetric Heavy Hydrocarbons	µg/L		500	NA	NA
<b>Surrogate</b>	<b>Unit</b>	<b>Acceptable Limits</b>			
Terphenyl	%	60-140	83	101	

Comments: RDL - Reported Detection Limit; G / S - Guideline / Standard: Refers to CCME FWAL-DCC

Guideline values are for general reference only. The guidelines provided may or may not be relevant for the intended use. Refer directly to the applicable standard for regulatory interpretation.

9381766-9381776 The C6-C10 fraction is calculated using Toluene response factor.

The C10 - C16, C16 - C34, and C34 - C50 fractions are calculated using the average response factor for n-C10, n-C16, and nC34.

Gravimetric Heavy Hydrocarbons are not included in the Total C16 - C50 and are only determined if the chromatogram of the C34 - C50 Hydrocarbons indicated that hydrocarbons &gt;C50 are present.

The chromatogram has returned to baseline by the retention time of nC50.

Total C6-C50 results are corrected for BTEX contributions.

This method complies with the Reference Method for the CWS PHC and is validated for use in the laboratory.

nC6 and nC10 response factors are within 30% of Toluene response factor.

nC10, nC16 and nC34 response factors are within 10% of their average.

C50 response factor is within 70% of nC10 + nC16 nC34 average.

Linearity is within 15%.

Extraction and holding times were met for this sample.

Fractions 1-4 are quantified with the contribution of PAHs. Under Ontario Regulation 153/04, results are considered valid without determining the PAH contribution if not requested by the client.

NA = Not Applicable

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**AGAT**

Laboratories

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## Polycyclic Aromatic Hydrocarbons in Water - (PAH)

DATE RECEIVED: 2018-07-06

DATE REPORTED: 2018-07-17

Parameter	Unit	SAMPLE DESCRIPTION:		ERK-4	ERK-5
		SAMPLE TYPE:		Water	Water
		G / S	DATE SAMPLED:	2018-07-01	2018-07-01
				9381766	9381776
1-Methylnaphthalene	ug/L		0.01	0.15	0.59
2-Methylnaphthalene	ug/L		0.01	0.20	0.80
Acenaphthene	ug/L		0.01	<0.01	<0.01
Acenaphthylene	ug/L		0.01	<0.01	<0.01
Acridine	ug/L		0.01	<0.01	<0.01
Anthracene	ug/L		0.012	<0.012	<0.012
Benzo(a)anthracene	ug/L		0.018	<0.018	<0.018
Benzo(a)pyrene	ug/L		0.010	<0.010	<0.010
Benzo(b)fluoranthene	ug/L		0.01	<0.01	<0.01
Benzo(e)pyrene	ug/L		0.01	<0.01	<0.01
Benzo(ghi)perylene	ug/L		0.01	<0.01	<0.01
Benzo(k)fluoranthene	ug/L		0.01	<0.01	<0.01
Chrysene	ug/L		0.01	<0.01	<0.01
Dibenzo(a,h)anthracene	ug/L		0.01	<0.01	<0.01
Fluoranthene	ug/L		0.01	<0.01	<0.01
Fluorene	ug/L		0.01	0.05	0.09
Indeno(1,2,3-cd)pyrene	ug/L		0.01	<0.01	<0.01
Naphthalene	ug/L		0.01	<0.01	<0.01
Perylene	ug/L		0.01	<0.01	<0.01
Phenanthrene	ug/L		0.01	<0.01	<0.01
Pyrene	ug/L		0.01	<0.01	<0.01
Quinoline	ug/L		0.01	<0.01	<0.01
Surrogate	Unit	Acceptable Limits			
Nitrobenzene-d5	%	50-140	93	91	
2-Fluorobiphenyl	%	50-140	76	82	
Terphenyl-d14	%	50-140	88	52	

Comments: RDL - Reported Detection Limit; G / S - Guideline / Standard: Refers to CCME FWAL-DCC

Guideline values are for general reference only. The guidelines provided may or may not be relevant for the intended use. Refer directly to the applicable standard for regulatory interpretation.

9381766-9381776 Benzo(b)fluoranthene may include contributions from benzo(j)fluoranthene, if also present in the sample.

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## Inorganic Chemistry (Water)

DATE RECEIVED: 2018-07-06

DATE REPORTED: 2018-07-17

		SAMPLE DESCRIPTION: ERK-1	
		SAMPLE TYPE: Water	
		DATE SAMPLED: 2018-07-01	
Parameter	Unit	G / S	RDL
BOD (5)	mg/L	5	70
pH	pH Units	NA	9.44
Total Suspended Solids	mg/L	10	221
Chemical Oxygen Demand	mg/L	50	456

Comments: RDL - Reported Detection Limit; G / S - Guideline / Standard

9381778 The RDL was raised for COD to reflect dilution of the sample in order to keep the analytes within a valid calibration range of the instruments.



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## Inorganic Chemistry (Water)

DATE RECEIVED: 2018-07-06

DATE REPORTED: 2018-07-17

Parameter	Unit	SAMPLE DESCRIPTION:		ERK-4	ERK-5
		SAMPLE TYPE:	DATE SAMPLED:	Water	Water
				2018-07-01	2018-07-01
Parameter	Unit	G / S	RDL	9381766	RDL
BOD (5)	mg/L		5	<5	<5
Electrical Conductivity	µS/cm		2	1040	2
pH	pH Units		NA	7.77	7.77
Total Suspended Solids	mg/L		10	<10	10
Total Hardness (as CaCO <sub>3</sub> )	mg/L		0.5	229	0.5
Alkalinity (as CaCO <sub>3</sub> )	mg/L		5	87	5
Chloride	mg/L		0.50	158	0.20
Nitrate as N	mg/L		0.25	<0.25	0.10
Nitrite as N	mg/L		0.25	<0.25	0.10
Sulphate	mg/L		0.50	159	0.20
Ammonia as N	mg/L		0.02	<0.02	<0.02
Chemical Oxygen Demand	mg/L		5	24	5
Phenols	µg/L		1	<1	1
Calcium	mg/L		0.10	58.1	0.05
Magnesium	mg/L		0.10	20.3	0.05
Sodium	mg/L		0.10	114	0.05
Potassium	mg/L		0.10	10.0	0.05
Total Arsenic	mg/L		0.015	<0.015	0.015
Total Cadmium	mg/L		0.010	<0.010	0.010
Total Chromium	mg/L		0.015	<0.015	0.015
Total Copper	mg/L		0.015	<0.015	0.015
Total Iron	mg/L		0.05	0.28	0.05
Total Lead	µg/L		0.5	<0.5	0.5
Total Mercury	mg/L		0.0002	<0.0002	0.0002
Total Nickel	mg/L		0.015	<0.015	<0.015

Comments: RDL - Reported Detection Limit; G / S - Guideline / Standard

9381766-9381776 Elevated RDLs indicate the degree of sample dilutions prior to analyses to keep analytes within the calibration range, reduce matrix interference and to avoid contaminating the instruments.

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## Quality Assurance

CLIENT NAME: DEFENCE CONSTRUCTION CANADA

AGAT WORK ORDER: 18T358833

PROJECT: Eureka Water License

ATTENTION TO: Cameron Chadwick

SAMPLING SITE:

SAMPLED BY:

### Microbiology Analysis

RPT Date: Jul 17, 2018			DUPLICATE			Method Blank	REFERENCE MATERIAL		METHOD BLANK SPIKE			MATRIX SPIKE			
PARAMETER	Batch	Sample Id	Dup #1	Dup #2	RPD		Measured Value	Acceptable Limits		Recovery	Acceptable Limits		Recovery	Acceptable Limits	
			Lower	Upper	Lower		Recovery	Lower	Upper	Lower	Upper	Lower	Recovery	Lower	Upper

**Microbiological Analysis (water)**

Fecal Coliform 9382031 ND ND NA < 1

Comments: ND – Not detected; NA - % RPD Not Applicable

**Certified By:**

*Nivine Basily*



## Quality Assurance

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### Trace Organics Analysis

RPT Date: Jul 17, 2018			DUPLICATE			Method Blank	REFERENCE MATERIAL			METHOD BLANK SPIKE			MATRIX SPIKE			
PARAMETER	Batch	Sample Id	Dup #1	Dup #2	RPD		Measured Value	Acceptable Limits		Recovery	Acceptable Limits		Recovery	Acceptable Limits		
								Lower	Upper		Lower	Upper		Lower	Upper	
<b>PHCs F1 - F4 (Water)</b>																
Benzene	9384209		< 0.20	< 0.20	NA	< 0.20	85%	50%	140%	89%	60%	130%	80%	50%	140%	
Toluene	9384209		< 0.20	< 0.20	NA	< 0.20	85%	50%	140%	82%	60%	130%	84%	50%	140%	
Ethylbenzene	9384209		< 0.10	< 0.10	NA	< 0.10	83%	50%	140%	82%	60%	130%	82%	50%	140%	
Xylene Mixture	9384209		< 0.20	< 0.20	NA	< 0.20	86%	50%	140%	84%	60%	130%	84%	50%	140%	
F1 (C6 - C10)	9384209		< 25	< 25	NA	< 25	90%	60%	140%	94%	60%	140%	82%	60%	140%	
F2 (C10 to C16)	TW		< 100	< 100	NA	< 100	103%	60%	140%	78%	60%	140%	75%	60%	140%	
F3 (C16 to C34)	TW		< 100	< 100	NA	< 100	102%	60%	140%	100%	60%	140%	106%	60%	140%	
F4 (C34 to C50)	TW		< 100	< 100	NA	< 100	86%	60%	140%	100%	60%	140%	101%	60%	140%	
<b>Oil and Grease (+Total) - water</b>																
Oil and Grease (animal/vegetable)	TW		< 0.5	< 0.5	NA	< 0.5	NA	70%	130%	101%	70%	130%	104%	70%	130%	
Oil and Grease (mineral) in water	TW		< 0.5	< 0.5	NA	< 0.5	NA	70%	130%	83%	70%	130%	84%	70%	130%	
Oil and Grease (Total) in water	TW		< 0.5	< 0.5	NA	< 0.5	NA	70%	130%	92%	70%	130%	94%	70%	130%	
<b>Polycyclic Aromatic Hydrocarbons in Water - (PAH)</b>																
1-Methylnaphthalene	1	9385494	< 0.01	< 0.01	NA	< 0.01	101%	50%	140%	96%	50%	140%	84%	50%	140%	
2-Methylnaphthalene	1	9385494	< 0.01	< 0.01	NA	< 0.01	99%	50%	140%	95%	50%	140%	83%	50%	140%	
Acenaphthene	1	9385494	< 0.01	< 0.01	NA	< 0.01	99%	50%	140%	95%	50%	140%	83%	50%	140%	
Acenaphthylene	1	9385494	< 0.01	< 0.01	NA	< 0.01	91%	50%	140%	94%	50%	140%	81%	50%	140%	
Acridine	1	9385494	< 0.01	< 0.01	NA	< 0.01	77%	50%	140%	98%	50%	140%	85%	50%	140%	
Anthracene	1	9385494	< 0.012	< 0.012	NA	< 0.012	89%	50%	140%	90%	50%	140%	79%	50%	140%	
Benzo(a)anthracene	1	9385494	< 0.018	< 0.018	NA	< 0.018	92%	50%	140%	91%	50%	140%	80%	50%	140%	
Benzo(a)pyrene	1	9385494	< 0.010	< 0.010	NA	< 0.010	81%	50%	140%	81%	50%	140%	71%	50%	140%	
Benzo(b)fluoranthene	1	9385494	< 0.01	< 0.01	NA	< 0.01	99%	50%	140%	88%	50%	140%	78%	50%	140%	
Benzo(e)pyrene	1	9385494	< 0.01	< 0.01	NA	< 0.01	92%	50%	140%	81%	50%	140%	72%	50%	140%	
Benzo(ghi)perylene	1	9385494	< 0.01	< 0.01	NA	< 0.01	76%	50%	140%	72%	50%	140%	62%	50%	140%	
Benzo(k)fluoranthene	1	9385494	< 0.01	< 0.01	NA	< 0.01	78%	50%	140%	76%	50%	140%	63%	50%	140%	
Chrysene	1	9385494	< 0.01	< 0.01	NA	< 0.01	94%	50%	140%	86%	50%	140%	78%	50%	140%	
Dibenzo(a,h)anthracene	1	9385494	< 0.01	< 0.01	NA	< 0.01	72%	50%	140%	78%	50%	140%	66%	50%	140%	
Fluoranthene	1	9385494	< 0.01	< 0.01	NA	< 0.01	92%	50%	140%	96%	50%	140%	83%	50%	140%	
Fluorene	1	9385494	< 0.01	< 0.01	NA	< 0.01	94%	50%	140%	96%	50%	140%	84%	50%	140%	
Indeno(1,2,3-cd)pyrene	1	9385494	< 0.01	< 0.01	NA	< 0.01	66%	50%	140%	82%	50%	140%	68%	50%	140%	
Naphthalene	1	9385494	< 0.01	< 0.01	NA	< 0.01	100%	50%	140%	92%	50%	140%	81%	50%	140%	
Perylene	1	9385494	< 0.01	< 0.01	NA	< 0.01	89%	50%	140%	81%	50%	140%	72%	50%	140%	
Phenanthrene	1	9385494	< 0.01	< 0.01	NA	< 0.01	96%	50%	140%	91%	50%	140%	79%	50%	140%	
Pyrene	1	9385494	< 0.01	< 0.01	NA	< 0.01	91%	50%	140%	96%	50%	140%	84%	50%	140%	
Quinoline	1	9385494	< 0.01	< 0.01	NA	< 0.01	96%	50%	140%	91%	50%	140%	80%	50%	140%	

Comments: If Matrix spike value is NA, the spiked analyte concentration was lower than that of the matrix contribution.

If RPD value is NA, the results of the duplicates are less than 5x the RDL and the RPD will not be calculated.



## Quality Assurance

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### Trace Organics Analysis (Continued)

RPT Date: Jul 17, 2018			DUPLICATE			Method Blank	REFERENCE MATERIAL		METHOD BLANK SPIKE			MATRIX SPIKE			
PARAMETER	Batch	Sample Id	Dup #1	Dup #2	RPD		Measured Value	Acceptable Limits	Recovery	Acceptable Limits		Recovery	Acceptable Limits		
							Lower	Upper		Lower	Upper		Lower	Upper	

**Certified By:** 



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### Water Analysis

RPT Date: Jul 17, 2018			DUPLICATE			Method Blank	REFERENCE MATERIAL		METHOD BLANK SPIKE			MATRIX SPIKE				
PARAMETER	Batch	Sample Id	Dup #1	Dup #2	RPD		Measured Value	Acceptable Limits		Recovery	Acceptable Limits		Recovery	Acceptable Limits		
								Lower	Upper			Lower		Recovery	Lower	Upper

#### Inorganic Chemistry (Water)

BOD (5)	9381766	9381766	<5	<5	NA	< 5	100%	75%	125%						
Electrical Conductivity	9376913		166	170	2.4%	< 2	102%	80%	120%						
pH	9376913		7.80	7.77	0.4%	NA	100%	90%	110%						
Total Suspended Solids	9378610		1230	1260	2.4%	< 10	102%	80%	120%						
Alkalinity (as CaCO <sub>3</sub> )	9376913		71	71	0.0%	< 5	106%	80%	120%						
Chloride	9381766	9381766	158	161	1.9%	< 0.10	97%	90%	110%	102%	90%	110%	95%	80%	120%
Nitrate as N	9381766	9381766	<0.25	<0.25	NA	< 0.05	101%	90%	110%	110%	90%	110%	111%	80%	120%
Nitrite as N	9381766	9381766	<0.25	<0.25	NA	< 0.05	NA	90%	110%	95%	90%	110%	115%	80%	120%
Sulphate	9381766	9381766	159	163	2.5%	< 0.10	104%	90%	110%	102%	90%	110%	96%	80%	120%
Ammonia as N	9381766	9381766	<0.02	<0.02	NA	< 0.02	104%	90%	110%	103%	90%	110%	103%	80%	120%
Chemical Oxygen Demand	9383252		26	28	7.4%	< 5	92%	80%	120%	95%	90%	110%	106%	70%	130%
Phenols	9381766	9381766	<1	<1	NA	< 1	102%	90%	110%	99%	90%	110%	98%	80%	120%
Calcium	9383573		67.6	67.1	0.7%	< 0.05	101%	90%	110%	101%	90%	110%	103%	70%	130%
Magnesium	9383573		22.1	22.2	0.5%	< 0.05	99%	90%	110%	99%	90%	110%	102%	70%	130%
Sodium	9383573		4.10	4.12	0.5%	< 0.05	92%	90%	110%	92%	90%	110%	93%	70%	130%
Potassium	9383573		1.83	1.88	2.7%	< 0.05	96%	90%	110%	96%	90%	110%	95%	70%	130%
Total Arsenic	9387771		<0.015	<0.015	NA	< 0.015	103%	90%	110%	97%	80%	120%	101%	70%	130%
Total Cadmium	9387771		<0.010	<0.010	NA	< 0.010	97%	90%	110%	99%	80%	120%	114%	70%	130%
Total Chromium	9387771		<0.015	<0.015	NA	< 0.015	106%	90%	110%	104%	80%	120%	103%	70%	130%
Total Copper	9387771		0.206	0.217	5.2%	< 0.015	105%	90%	110%	105%	80%	120%	100%	70%	130%
Total Iron	9387771		0.16	0.18	NA	< 0.05	110%	90%	110%	112%	80%	120%	110%	70%	130%
Total Lead	9387771		1.8	1.7	NA	< 0.5	105%	90%	110%	103%	80%	120%	100%	70%	130%
Total Mercury	9382031		<0.0002	<0.0002	NA	< 0.0002	105%	90%	110%	104%	90%	110%	97%	80%	120%
Total Nickel	9387771		0.025	0.026	NA	< 0.015	106%	90%	110%	103%	80%	120%	101%	70%	130%

Comments: NA signifies Not Applicable.

Duplicate Qualifier: As the measured result approaches the Reporting Limit (RL), the uncertainty associated with the value increases dramatically, thus duplicate acceptance limits apply only where the average of the two duplicates is greater than five times the RL.

**Certified By:**





## Method Summary

CLIENT NAME: DEFENCE CONSTRUCTION CANADA

AGAT WORK ORDER: 18T358833

PROJECT: Eureka Water License

ATTENTION TO: Cameron Chadwick

SAMPLING SITE:

SAMPLED BY:

PARAMETER	AGAT S.O.P	LITERATURE REFERENCE	ANALYTICAL TECHNIQUE
<b>Microbiology Analysis</b>			
Fecal Coliform	MIC-93-7000	SM 9222 D	Membrane Filtration
<b>Trace Organics Analysis</b>			
Oil and Grease (animal/vegetable)	VOL-91- 5011	SM 5520 & EPA SW846 3510C & EPA 1664	GRAVIMETRIC
Oil and Grease (mineral) in water	VOL-91- 5011	SM 5520 & EPA SW846 3510C & EPA 1664	GRAVIMETRIC
Oil and Grease (Total) in water	VOL-91- 5011	SM 5520 & EPA SW846 3510C & EPA 1664	GRAVIMETRIC
Benzene	VOL-91-5010	MOE PHC-E3421	(P&T)GC/FID
Toluene	VOL-91-5010	MOE PHC-E3421	(P&T)GC/FID
Ethylbenzene	VOL-91-5010	MOE PHC-E3421	(P&T)GC/FID
Xylene Mixture	VOL-91-5010	MOE PHC-E3421	(P&T)GC/FID
F1 (C6 - C10)	VOL-91- 5010	MOE PHC-E3421	(P&T)GC/FID
F1 (C6 to C10) minus BTEX	VOL-91-5010	MOE PHC-E3421	(P&T)GC/FID
F2 (C10 to C16)	VOL-91-5010	MOE PHC-E3421	GC/FID
F3 (C16 to C34)	VOL-91-5010	MOE PHC-E3421	GC/FID
F4 (C34 to C50)	VOL-91-5010	MOE PHC-E3421	GC/FID
Gravimetric Heavy Hydrocarbons	VOL-91-5010	MOE PHC-E3421	BALANCE
Terphenyl	VOL-91-5010		GC/FID
1-Methylnaphthalene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
2-Methylnaphthalene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Acenaphthene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Acenaphthylene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Acridine	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Anthracene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Benzo(a)anthracene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Benzo(a)pyrene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Benzo(b)fluoranthene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Benzo(e)pyrene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Benzo(ghi)perylene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Benzo(k)fluoranthene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Chrysene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Dibenzo(a,h)anthracene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Fluoranthene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Fluorene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Indeno(1,2,3-cd)pyrene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Naphthalene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Perylene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Phenanthrene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Pyrene	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Quinoline	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Nitrobenzene-d5	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
2-Fluorobiphenyl	ORG-120-5104	EPA SW846/3510/8270C	GC/MS
Terphenyl-d14	ORG-120-5104	EPA SW846/3510/8270C	GC/MS



## Method Summary

CLIENT NAME: DEFENCE CONSTRUCTION CANADA

AGAT WORK ORDER: 18T358833

PROJECT: Eureka Water License

ATTENTION TO: Cameron Chadwick

SAMPLING SITE:

SAMPLED BY:

PARAMETER	AGAT S.O.P	LITERATURE REFERENCE	ANALYTICAL TECHNIQUE
<b>Water Analysis</b>			
BOD (5)	INOR-93-6006	SM 5210 B	DO METER
pH	INOR-93-6000	SM 4500-H+ B	PC TITRATE
Total Suspended Solids	INOR-93-6028	SM 2540 D	BALANCE
Chemical Oxygen Demand	INOR-93-6042	SM 5220 D	SPECTROPHOTOMETER
Electrical Conductivity	INOR-93-6000	SM 2510 B	PC TITRATE
Total Hardness (as CaCO <sub>3</sub> )	MET-93-6105	EPA SW-846 6010C & 200.7	CALCULATION
Alkalinity (as CaCO <sub>3</sub> )	INOR-93-6000	SM 2320 B	PC TITRATE
Chloride	INOR-93-6004	SM 4110 B	ION CHROMATOGRAPH
Nitrate as N	INOR-93-6004	SM 4110 B	ION CHROMATOGRAPH
Nitrite as N	INOR-93-6004	SM 4110 B	ION CHROMATOGRAPH
Sulphate	INOR-93-6004	SM 4110 B	ION CHROMATOGRAPH
Ammonia as N	INOR-93-6059	QuikChem 10-107-06-1-J & SM 4500 NH3-F	LACHAT FIA
Phenols	INOR-93-6050	MOE ROPHEN-E 3179 & SM 5530 D	TECHNICON AUTO ANALYZER
Calcium	MET-93-6105	EPA SW-846 6010C & 200.7	ICP/OES
Magnesium	MET-93-6105	EPA SW-846 6010C & 200.7	ICP/OES
Sodium	MET-93-6105	EPA SW-846 6010C & 200.7	ICP/OES
Potassium	MET-93-6105	EPA SW-846 6010C & 200.7	ICP/OES
Total Arsenic	MET-93-6103	EPA SW-846 3010A & 6020A	ICP-MS
Total Cadmium	MET-93-6103	EPA SW-846 3010A & 6020A	ICP-MS
Total Chromium	MET-93-6103	EPA SW-846 3010A & 6020A	ICP-MS
Total Copper	MET-93-6103	EPA SW-846 3010A & 6020A	ICP-MS
Total Iron	MET-93-6103	EPA SW-846 3010A & 6020A	ICP-MS
Total Lead	MET-93-6003	EPA SW-846 3010A & 6020A	ICP/MS
Total Mercury	MET-93-6100	EPA SW-846 7470 & 245.1	CVAAS
Total Nickel	MET-93-6103	EPA SW-846 3010A & 6020A	ICP-MS





## Sample Temperature Log

Client: DND

# of Coolers: 6 large + 1 med blue

### Arrival Temperatures - Branch/Driver

Cooler #1: 4.7 / 4.9 / 4.8

Cooler #2: 8.8 / 9.0 / 9.1

Cooler #3: 7.4 / 7.5 / 7.6

Cooler #4: 6.9 / 7.0 / 6.8

Cooler #5: 7.8 / 7.6 / 8.0

Cooler #6: 5.8 / 6.0 / 6.2

Cooler #7: 8.1 / 8.4 / 8.5

Cooler #8:   /  /  

Cooler #9:   /  /  

Cooler #10:   /  /  

IR Gun ID:   

Taken By: Arneet Minhas

Date (yyyy/mm/dd): 2018/07/06 Time: 9 : 48 AM / PM

COC# or Work Order #:   

# of Submissions:   

### Arrival Temperatures - Laboratory

Cooler #1:   /  /  

Cooler #2:   /  /  

Cooler #3:   /  /  

Cooler #4:   /  /  

Cooler #5:   /  /  

Cooler #6:   /  /  

Cooler #7:   /  /  

Cooler #8:   /  /  

Cooler #9:   /  /  

Cooler #10:   /  /  

IR Gun ID:   

Taken By:   

Date

(yyyy/mm/dd):    Time:   ;    AM / PM

Instructions for use of this form: 1) complete all fields of info including total # of coolers and # of submissions rec'd, 2) photocopy and place in each submission prior to giving a WO#, 3) Proceed as normal, write the WO# and scan ( please make sure to scan along with the COC)