

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\WX050120\
 Data File : VX016017.D
 Acq On : 01 May 2020 17:25
 Operator : JC/SP
 Sample : L2469-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SA-PZ187I-20200429

Manual Integrations
 APPROVED

MMDadoda
 5/4/2020 2:19:59 PM

Quant Time: May 04 04:53:42 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042820W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 28 17:30:11 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	276106	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	451083	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	436267	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	225823	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	176397	48.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.72%	
35) Dibromofluoromethane	5.47	113	132822	47.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.40%	
50) Toluene-d8	8.70	98	550458	52.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.64%	
62) 4-Bromofluorobenzene	11.12	95	223334	54.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.24%	
Target Compounds						
6) Chloroethane	1.72	64	7301	3.228	ug/l	98
12) 1,1-Dichloroethene	2.36	96	2109	0.798	ug/l	98
16) Acetone	2.42	43	2415	1.564	ug/l	97
24) 1,1-Dichloroethane	3.67	63	59328	11.372	ug/l	97
85) sec-Butylbenzene	11.93	105	5653	0.376	ug/l	84
87) 1,3-Dichlorobenzene	12.01	146	3194	0.443	ug/l	95
88) 1,4-Dichlorobenzene	12.08	146	9723m	1.334	ug/l	
91) 1,2-Dichlorobenzene	12.38	146	4734	0.687	ug/l	99
93) 1,2,4-Trichlorobenzene	13.63	180	7031	1.358	ug/l	92

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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