

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX050120\
 Data File : VX016015.D
 Acq On : 01 May 2020 16:39
 Operator : JC/SP
 Sample : L2469-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SA-MW128I-20200428

Quant Time: May 02 07:52:08 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	275099	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	449845	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	439309	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	229715	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	177370	49.31	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.62%	
35) Dibromofluoromethane	5.47	113	135211	48.18	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.36%	
50) Toluene-d8	8.70	98	552174	52.63	ug/l	0.00
Spiked Amount	50.000		Recovery	=	105.26%	
62) 4-Bromofluorobenzene	11.12	95	220269	53.53	ug/l	0.00
Spiked Amount	50.000		Recovery	=	107.06%	
Target Compounds						
6) Chloroethane	1.72	64	4512	2.002	ug/l	99
12) 1,1-Dichloroethene	2.36	96	2289	0.870	ug/l	86
16) Acetone	2.42	43	3553	2.310	ug/l	98
19) Methyl tert-butyl Ether	3.17	73	9756	1.025	ug/l	92
24) 1,1-Dichloroethane	3.67	63	84667	16.289	ug/l	99
65) Chlorobenzene	10.12	112	3662	0.409	ug/l	94

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

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