

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112420\
 Data File : VX019639.D
 Acq On : 25 Nov 2020 05:05
 Operator : JC/MD
 Sample : L4835-21
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MH-121.5T-20201119

Quant Time: Nov 25 07:11:11 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
 Quant Title : SW846 8260
 QLast Update : Sat Nov 21 01:57:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	303263	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	498469	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.09	117	456101	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	182298	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	6.03	65	187843	47.77	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	95.54%
35) Dibromofluoromethane	5.46	113	148520	47.47	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	94.94%
50) Toluene-d8	8.70	98	596041	49.02	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	98.04%
62) 4-Bromofluorobenzene	11.12	95	206377	47.18	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	94.36%

Target Compounds

					Qvalue
16) Acetone	2.42	43	3983	2.619	ug/l 97
27) cis-1,2-Dichloroethene	4.56	96	21810	5.566	ug/l 93
30) Chloroform	5.17	83	16028	2.568	ug/l 94
44) Trichloroethene	7.18	130	9417	2.504	ug/l 99
47) Bromodichloromethane	7.87	83	4489	0.956	ug/l 96
60) Dibromochloromethane	9.56	129	1976	0.564	ug/l 95

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

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