

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112120\
 Data File : VX019531.D
 Acq On : 21 Nov 2020 06:41
 Operator : JC/MD
 Sample : L4826-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB-2020-WS-000-04

Quant Time: Nov 21 07:56:15 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	327619	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	527688	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	478558	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	208215	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	6.03	65	193909	45.65	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	91.30%		
35) Dibromofluoromethane	5.46	113	157022	47.41	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	94.82%		
50) Toluene-d8	8.70	98	630042	48.95	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	97.90%		
62) 4-Bromofluorobenzene	11.12	95	229026	49.46	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	98.92%		

Target Compounds

					Qvalue
16) Acetone	2.42	43	37042	22.549	ug/l 99
18) Methyl Acetate	2.75	43	2869	1.338	ug/l # 86
20) Methylene Chloride	2.83	84	18526	4.512	ug/l 97
43) Isopropyl Acetate	6.42	43	34446	4.412	ug/l 99
64) Tetrachloroethene	9.32	164	3773	1.003	ug/l 98

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

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