

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112120\
 Data File : VX019519.D
 Acq On : 21 Nov 2020 02:00
 Operator : JC/MD
 Sample : L4779-02 10X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1641

Quant Time: Nov 21 06:10:00 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	5.62	168	259209	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	424787	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	366916	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	167057	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	6.03	65	158624	47.20	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	94.40%		
35) Dibromofluoromethane	5.46	113	126649	47.50	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	95.00%		
50) Toluene-d8	8.70	98	497365	48.00	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	96.00%		
62) 4-Bromofluorobenzene	11.12	95	176789	47.42	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	94.84%		

Target Compounds

					Qvalue
16) Acetone	2.44	43	26803	20.622 ug/l #	1
18) Methyl Acetate	2.77	43	3360	1.675 ug/l	98
25) 2-Butanone	4.66	43	6658	3.410 ug/l #	86
30) Chloroform	5.17	83	33946	6.364 ug/l	99
51) 4-Methyl-2-Pentanone	8.63	43	4990	1.362 ug/l	93
59) 2-Hexanone	9.48	43	1732	0.623 ug/l	92
95) Naphthalene	13.82	128	15711	3.398 ug/l #	95

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

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