

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112020\
 Data File : VX019483.D
 Acq On : 19 Nov 2020 22:00
 Operator : JC/MD
 Sample : L4801-02 5X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EFF-WW

Quant Time: Nov 20 03:04:41 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA X\METHOD\624X112020W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Nov 20 01:06:02 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	4.98	128	44444	30.00	ug/l	0.00
28) 1,4-Difluorobenzene	6.82	114	237552	30.00	ug/l	0.00
57) Chlorobenzene-d5	10.10	117	224430	30.00	ug/l	0.00

System Monitoring Compounds

27) 1,2-Dichloroethane-d4	6.03	65	97545	30.54	ug/l	0.00
Spiked Amount	30.000	Range	91 - 110	Recovery	=	101.80%
60) 4-Bromofluorobenzene	11.12	95	108050	29.52	ug/l	0.00
Spiked Amount	30.000	Range	63 - 112	Recovery	=	98.40%
63) Toluene-d8	8.70	98	290917	29.08	ug/l	0.00
Spiked Amount	30.000	Range	91 - 112	Recovery	=	96.93%

Target Compounds

					Ovalue
12) Methyl Acetate	2.75	43	17533	5.943	ug/l 95
15) Acetone	2.42	58	359476	860.286	ug/l 98
18) Methylene Chloride	2.83	84	1200	0.403	ug/l 92
20) Diisopropyl ether	3.82	45	3344	0.421	ug/l # 78
25) Chloroform	5.17	83	6706	1.321	ug/l 88
30) 2-Butanone	4.64	43	44326	24.041	ug/l 100
67) m/p-Xylenes	10.34	106	3056	0.593	ug/l 93
68) o-Xylene	10.68	106	1141	0.234	ug/l 96

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

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