

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102320\
 Data File : VX019104.D
 Acq On : 24 Oct 2020 04:06
 Operator : JC/SP
 Sample : L4417-18
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-10D

Quant Time: Oct 26 06:23:19 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 22 17:51:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	261225	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	452235	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	428672	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	223619	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	159888	51.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.20%	
35) Dibromofluoromethane	5.46	113	140316	49.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.02%	
50) Toluene-d8	8.70	98	561342	51.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.76%	
62) 4-Bromofluorobenzene	11.12	95	201452	51.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.66%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	3.01	59	12501	19.726	ug/l	99
16) Acetone	2.42	43	3488	3.273	ug/l	94
17) Carbon Disulfide	2.56	76	3974	0.542	ug/l #	90
19) Methyl tert-butyl Ether	3.17	73	29261	3.633	ug/l	97
67) Ethyl Benzene	10.24	91	4836	0.312	ug/l	91
68) m/p-Xylenes	10.34	106	2453	0.413	ug/l	93

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

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