

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031521\
 Data File : VX021224.D
 Acq On : 16 Mar 2021 01:31
 Operator : JC/MD
 Sample : M1647-22
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 COMP3-30788:89

Quant Time: Mar 16 04:42:33 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031521W.M
 Quant Title : SW846 8260
 QLast Update : Mon Mar 15 11:55:59 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.629	168	72509	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.829	114	134190	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	126598	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	58111	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.029	65	63247	51.15	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 102.30%	
35) Dibromofluoromethane	5.470	113	45957	49.78	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 99.56%	
50) ToLuene-d8	8.694	98	168540	50.82	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 101.64%	
62) 4-Bromofluorobenzene	11.118	95	67197	53.17	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 106.34%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	3.011	59	18262	84.80	ug/l	98
16) Acetone	2.423	43	43337	87.20	ug/l	89
17) Carbon Disulfide	2.552	76	1979	0.86	ug/l	96
18) Methyl Acetate	2.752	43	24116	20.04	ug/l #	88
25) 2-Butanone	4.635	43	7722	10.80	ug/l #	85
40) Benzene	6.117	78	62363	15.86	ug/l	98
51) 4-Methyl-2-Pentanone	8.618	43	1919	1.30	ug/l	87
52) ToLuene	8.765	92	21976	9.27	ug/l	100
67) Ethyl Benzene	10.235	91	4859	1.00	ug/l	99
68) m/p-Xylenes	10.341	106	8587	4.92	ug/l	93
69) o-Xylene	10.683	106	3935	2.32	ug/l	99
70) Styrene	10.694	104	2364	0.82	ug/l #	71
84) 1,2,4-Trimethylbenzene	11.789	105	3487	0.91	ug/l	83
95) Naphthalene	13.818	128	18007	5.55	ug/l	99

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

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