

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031521\
 Data File : VX021211.D
 Acq On : 15 Mar 2021 19:41
 Operator : JC/MD
 Sample : M1647-20
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 COMP4-41735-30780

Quant Time: Mar 16 03:05:41 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	70559	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.829	114	130433	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	123367	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	54872	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.029	65	61959	51.50	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 103.00%			
35) Dibromofluoromethane	5.470	113	45062	50.21	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 100.42%			
50) Toluene-d8	8.694	98	162311	50.35	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 100.70%			
62) 4-Bromofluorobenzene	11.118	95	64389	52.41	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 104.82%			
Target Compounds					Qvalue	
11) Tert butyl alcohol	3.011	59	3143	15.00	ug/l #	86
16) Acetone	2.422	43	564016	1166.23	ug/l	89
18) Methyl Acetate	2.752	43	2438	2.08	ug/l #	89
25) 2-Butanone	4.640	43	11157	16.04	ug/l	89
40) Benzene	6.117	78	22889	5.99	ug/l	97
51) 4-Methyl-2-Pentanone	8.617	43	3352	2.33	ug/l	89
52) Toluene	8.764	92	5119	2.22	ug/l	98
59) 2-Hexanone	9.470	43	1306	1.20	ug/l	77
67) Ethyl Benzene	10.229	91	1945	0.41	ug/l	100
68) m/p-Xylenes	10.341	106	4793	2.82	ug/l	90
69) o-Xylene	10.682	106	3757	2.28	ug/l	89
80) 1,3,5-Trimethylbenzene	11.488	105	3574	1.00	ug/l	98
84) 1,2,4-Trimethylbenzene	11.788	105	8528	2.37	ug/l	95
95) Naphthalene	13.818	128	4911	2.64	ug/l #	95

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

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