

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031521\
 Data File : VX021229.D
 Acq On : 16 Mar 2021 03:28
 Operator : JC/MD
 Sample : M1667-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DRUM-TOTE-COMP

Quant Time: Mar 16 04:44:30 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	5.635	168	67705	50.00	ug/l	0.01
34) 1,4-Difluorobenzene	6.829	114	121840	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	117752	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	54797	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.035	65	57779	50.05	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 100.10%			
35) Dibromofluoromethane	5.464	113	42613	50.83	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 101.66%			
50) Toluene-d8	8.694	98	152882	50.77	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 101.54%			
62) 4-Bromofluorobenzene	11.118	95	62646	54.59	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 109.18%			
Target Compounds						Qvalue
11) Tert butyl alcohol	3.011	59	2383	11.85	ug/l #	93
16) Acetone	2.423	43	13440	28.96	ug/l #	86
18) Methyl Acetate	2.752	43	243413	216.57	ug/l #	88
19) Methyl tert-butyl Ether	3.170	73	1708	0.62	ug/l	94
25) 2-Butanone	4.641	43	2422	3.63	ug/l #	87
31) Cyclohexane	5.547	56	5552	3.80	ug/l #	89
51) 4-Methyl-2-Pentanone	8.617	43	1656	1.23	ug/l #	85
52) Toluene	8.765	92	911	0.42	ug/l	90
68) m/p-Xylenes	10.341	106	742	0.46	ug/l	94
69) o-Xylene	10.682	106	686	0.44	ug/l	88
73) Isopropylbenzene	11.000	105	1089	0.26	ug/l	92
78) n-propylbenzene	11.341	91	1696	0.35	ug/l	90
84) 1,2,4-Trimethylbenzene	11.789	105	2009	0.56	ug/l	92

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

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