

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112421\
 Data File : VX025338.D
 Acq On : 24 Nov 2021 19:42
 Operator : JC/MD
 Sample : M4838-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 211123015-01-VOA

Quant Time: Nov 26 10:39:02 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 17 15:36:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	131258	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	207771	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	185696	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	85610	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	68792	49.721	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	99.440%		
35) Dibromofluoromethane	5.391	113	62005	48.638	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	97.280%		
50) Toluene-d8	8.653	98	237364	49.522	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	99.040%		
62) 4-Bromofluorobenzene	11.079	95	83615	46.112	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	92.220%		
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.959	59	925712	2795.927	ug/l	99
16) Acetone	2.374	43	3352	4.360	ug/l	95
29) Tetrahydrofuran	5.007	42	132364	192.330	ug/l	98
40) Benzene	6.031	78	3461	0.630	ug/l	95
49) 1,4-Dioxane	7.659	88	2860	84.075	ug/l	95
52) Toluene	8.720	92	1144	0.319	ug/l	90
65) Chlorobenzene	10.079	112	1528	0.404	ug/l	91
67) Ethyl Benzene	10.195	91	3143	0.481	ug/l	93
68) m/p-Xylenes	10.305	106	3190	1.194	ug/l	98
69) o-Xylene	10.646	106	2482	0.949	ug/l	85
73) Isopropylbenzene	10.963	105	4394	0.720	ug/l	94
84) 1,2,4-Trimethylbenzene	11.756	105	2445	0.482	ug/l	94

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

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