

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX120319\
 Data File : VX013739.D
 Acq On : 03 Dec 2019 20:51
 Operator : JC/SP
 Sample : K6084-06 20X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 30738

Quant Time: Dec 04 08:12:07 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 26 15:53:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	747531	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	1117022	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	1005464	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	484047	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.05	65	394268	45.44	ug/l	0.00
Spiked Amount			Recovery	=	90.88%	
35) Dibromofluoromethane	5.49	113	337781	48.91	ug/l	0.00
Spiked Amount			Recovery	=	97.82%	
50) Toluene-d8	8.71	98	1331239	49.48	ug/l	0.00
Spiked Amount			Recovery	=	98.96%	
62) 4-Bromofluorobenzene	11.14	95	472499	48.70	ug/l	0.00
Spiked Amount			Recovery	=	97.40%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
11) Tert butyl alcohol	3.02	59	252485	117.885	ug/l	98
16) Acetone	2.43	43	419802	97.993	ug/l	97
18) Methyl Acetate	2.76	43	54906	4.603	ug/l	98
19) Methyl tert-butyl Ether	3.18	73	267572	11.354	ug/l	99
25) 2-Butanone	4.66	43	152629	23.972	ug/l	98
30) Chloroform	5.20	83	48256	3.501	ug/l	96
40) Benzene	6.14	78	2103816	68.124	ug/l	99
43) Isopropyl Acetate	6.43	43	35641	1.903	ug/l #	91
49) 1,4-Dioxane	7.73	88	8357	41.031	ug/l #	56
52) Toluene	8.78	92	2316626	120.750	ug/l	100
67) Ethyl Benzene	10.25	91	209995	5.854	ug/l	99
68) m/p-Xylenes	10.35	106	579768	42.214	ug/l	99
69) o-Xylene	10.70	106	180581	13.482	ug/l	95
80) 1,3,5-Trimethylbenzene	11.50	105	98152	3.527	ug/l	97
84) 1,2,4-Trimethylbenzene	11.81	105	209455	7.416	ug/l	99

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

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