

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX041218\
 Data File : VX000808.D
 Acq On : 12 Apr 2018 17:48
 Operator : JC/MD
 Sample : J2236-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-103

Quant Time: Apr 13 07:31:34 2018
 Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X041018W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 10 15:49:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.68	168	274872	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	359491	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	322697	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	208500	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.08	65	156198	45.07	ug/l	0.00
Spiked Amount			Recovery	=	90.14%	
35) Dibromofluoromethane	5.51	113	129641	46.81	ug/l	0.00
Spiked Amount			Recovery	=	93.62%	
50) Toluene-d8	8.73	98	491840	47.69	ug/l	0.00
Spiked Amount			Recovery	=	95.38%	
62) 4-Bromofluorobenzene	11.15	95	190124	44.56	ug/l	0.00
Spiked Amount			Recovery	=	89.12%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.45	43	28823	24.12	ug/l	100
17) Carbon Disulfide	2.57	76	4670	0.61	ug/l	97
19) Methyl tert-butyl Ether	3.21	73	5868	0.70	ug/l	98
25) 2-Butanone	4.73	43	7483	3.79	ug/l	100
27) cis-1,2-Dichloroethene	4.61	96	1475	0.47	ug/l	88
30) Chloroform	5.22	83	6904	1.36	ug/l	91
40) Benzene	6.15	78	15087	1.37	ug/l	95
44) Trichloroethene	7.23	130	2944	0.86	ug/l	87
52) Toluene	8.79	92	4309	0.59	ug/l	99
64) Tetrachloroethene	9.34	164	5984	1.90	ug/l	98
67) Ethyl Benzene	10.26	91	7070	0.52	ug/l	98
68) m/p-Xylenes	10.36	106	3835	0.72	ug/l	97
69) o-Xylene	10.71	106	2406	0.47	ug/l	99
70) Styrene	10.72	104	4618	0.53	ug/l	95
84) 1,2,4-Trimethylbenzene	11.81	105	4207	0.36	ug/l	98
95) Naphthalene	13.84	128	126387	8.33	ug/l	100

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

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