

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX050818\
 Data File : VX001506.D
 Acq On : 08 May 2018 11:46
 Operator : JC/MD
 Sample : J2695-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CS-DRUM-2

Quant Time: May 08 14:12:52 2018
 Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X050518W.M
 Quant Title : SW846 8260
 QLast Update : Sun May 06 07:59:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	261937	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	358353	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	326607	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.09	152	196556	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.07	65	175130	46.52	ug/l	0.00
Spiked Amount			Recovery	=	93.04%	
35) Dibromofluoromethane	5.51	113	136120	42.70	ug/l	0.00
Spiked Amount			Recovery	=	85.40%	
50) Toluene-d8	8.72	98	514828	44.25	ug/l	0.00
Spiked Amount			Recovery	=	88.50%	
62) 4-Bromofluorobenzene	11.14	95	178244	39.91	ug/l	0.00
Spiked Amount			Recovery	=	79.82%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
11) Tert butyl alcohol	3.07	59	7322	9.22	ug/l	98
16) Acetone	2.45	43	509590	298.09	ug/l	99
17) Carbon Disulfide	2.57	76	10583	1.43	ug/l	98
25) 2-Butanone	4.71	43	218764	88.53	ug/l	98
30) Chloroform	5.22	83	15331	2.64	ug/l	97
40) Benzene	6.15	78	4628	0.35	ug/l	97
51) 4-Methyl-2-Pentanone	8.65	43	178879	35.09	ug/l	99
52) Toluene	8.79	92	10079	1.16	ug/l	99
59) 2-Hexanone	9.50	43	23889	6.12	ug/l	100
67) Ethyl Benzene	10.26	91	3879	0.25	ug/l	99
68) m/p-Xylenes	10.36	106	5769	0.93	ug/l	95
69) o-Xylene	10.70	106	2914	0.48	ug/l	98
84) 1,2,4-Trimethylbenzene	11.81	105	8093	0.70	ug/l	92
86) p-Isopropyltoluene	12.07	119	3460	0.28	ug/l #	71
95) Naphthalene	13.84	128	5300	0.36	ug/l	96

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

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