

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX041019\
 Data File : VX008842.D
 Acq On : 11 Apr 2019 06:34
 Operator : JC/SP
 Sample : K2344-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WOOD-2

Quant Time: Apr 11 09:06:59 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X040319W.M
 Quant Title : SW846 8260
 QLast Update : Thu Apr 11 08:05:56 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	187326	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	304767	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	266500	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	112642	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	125484	45.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.86%	
35) Dibromofluoromethane	5.49	113	91632	47.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.06%	
50) Toluene-d8	8.71	98	385119	49.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.28%	
62) 4-Bromofluorobenzene	11.13	95	129622	47.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.22%	
Target Compounds						
					Qvalue	
13) Acrolein	2.29	56	2596	8.185	ug/l	97
16) Acetone	2.44	43	18744	11.911	ug/l	100
18) Methyl Acetate	2.78	43	5124	1.343	ug/l	99
20) Methylene Chloride	2.85	84	58695	22.282	ug/l	97
25) 2-Butanone	4.68	43	3823	1.537	ug/l	92
43) Isopropyl Acetate	6.45	43	10017	1.438	ug/l	96
80) 1,3,5-Trimethylbenzene	11.51	105	8833	1.127	ug/l	99
84) 1,2,4-Trimethylbenzene	11.80	105	24141	3.043	ug/l	100

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

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