

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX041019\
 Data File : VX008841.D
 Acq On : 11 Apr 2019 06:11
 Operator : JC/SP
 Sample : K2344-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WOOD-1

Quant Time: Apr 11 09:04:49 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	5.67	168	192746	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	312520	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	274795	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	116821	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	129843	46.19	ug/l	0.00
Spiked Amount	50.000		Recovery	=	92.38%	
35) Dibromofluoromethane	5.50	113	94282	47.19	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.38%	
50) Toluene-d8	8.71	98	395852	49.75	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.50%	
62) 4-Bromofluorobenzene	11.13	95	134350	47.62	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.24%	
Target Compounds						
					Qvalue	
13) Acrolein	2.29	56	3222	9.873	ug/l	91
16) Acetone	2.44	43	22076	13.634	ug/l	99
18) Methyl Acetate	2.78	43	5464	1.391	ug/l	97
20) Methylene Chloride	2.85	84	342147	126.235	ug/l	97
25) 2-Butanone	4.69	43	4623	1.806	ug/l	96
43) Isopropyl Acetate	6.45	43	10864	1.521	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX041019\
Data File : VX008841.D
Acq On : 11 Apr 2019 06:11
Operator : JC/SP
Sample : K2344-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WOOD-1

Quant Time: Apr 11 09:04:49 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X040319W.M
Quant Title : SW846 8260
QLast Update : Thu Apr 11 08:05:56 2019
Response via : Initial Calibration

