

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX041019\
 Data File : VX008833.D
 Acq On : 11 Apr 2019 03:05
 Operator : JC/SP
 Sample : K2200-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EO-01-040919-C

Quant Time: Apr 11 08:31:52 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X040319W.M
 Quant Title : SW846 8260
 QLast Update : Thu Apr 04 04:47:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	192476	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	314421	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	279573	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	121746	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	129433	46.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.20%	
35) Dibromofluoromethane	5.50	113	94695	47.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.22%	
50) Toluene-d8	8.71	98	402090	50.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.46%	
62) 4-Bromofluorobenzene	11.13	95	134474	47.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.76%	
Target Compounds						
16) Acetone	2.44	43	7624	4.715	ug/l	99
20) Methylene Chloride	2.85	84	3767740	1392.053	ug/l	98
43) Isopropyl Acetate	6.45	43	10181	1.417	ug/l	99
95) Naphthalene	13.83	128	135567	13.531	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX041019\
Data File : VX008833.D
Acq On : 11 Apr 2019 03:05
Operator : JC/SP
Sample : K2200-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EO-01-040919-C

Quant Time: Apr 11 08:31:52 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X040319W.M
Quant Title : SW846 8260
QLast Update : Thu Apr 04 04:47:41 2019
Response via : Initial Calibration

