

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX010919\
 Data File : VX006962.D
 Acq On : 09 Jan 2019 13:53
 Operator : JC/MD
 Sample : K1082-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SP-1-2

Quant Time: Jan 10 01:27:48 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 08 14:06:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	667623	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	1021809	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	954088	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	471009	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	351609	50.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.78%	
35) Dibromofluoromethane	5.51	113	313449	47.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.80%	
50) Toluene-d8	8.71	98	1216613	50.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.58%	
62) 4-Bromofluorobenzene	11.13	95	432264	51.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.54%	
Target Compounds						
8) Diethyl Ether	2.19	74	8757	1.984	ug/l	95
16) Acetone	2.45	43	29759	8.627	ug/l	99
20) Methylene Chloride	2.86	84	445472	68.106	ug/l	94
43) Isopropyl Acetate	6.46	43	32663	2.174	ug/l	98
95) Naphthalene	13.83	128	461815	13.609	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX010919\
Data File : VX006962.D
Acq On : 09 Jan 2019 13:53
Operator : JC/MD
Sample : K1082-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SP-1-2

Quant Time: Jan 10 01:27:48 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260
QLast Update : Tue Jan 08 14:06:59 2019
Response via : Initial Calibration

