

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX070219\
 Data File : VX010621.D
 Acq On : 02 Jul 2019 16:47
 Operator : JC/SP
 Sample : K3636-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SP19-JMW0961

Quant Time: Jul 03 03:50:43 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 20 03:31:56 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	243555	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	372033	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	342663	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	158818	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	118127	51.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.46%	
35) Dibromofluoromethane	5.50	113	114671	50.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.76%	
50) Toluene-d8	8.71	98	441951	50.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.50%	
62) 4-Bromofluorobenzene	11.13	95	149838	47.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.32%	
Target Compounds						
8) Diethyl Ether	2.19	74	29879	24.337	ug/l	98
24) 1,1-Dichloroethane	3.70	63	4153	1.147	ug/l	96
29) Tetrahydrofuran	5.13	42	3386	3.256	ug/l #	89
40) Benzene	6.14	78	10081	1.095	ug/l #	87

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070219\
Data File : VX010621.D
Acq On : 02 Jul 2019 16:47
Operator : JC/SP
Sample : K3636-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SP19-JMW0961

Quant Time: Jul 03 03:50:43 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260
QLast Update : Thu Jun 20 03:31:56 2019
Response via : Initial Calibration

