

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX050919\
 Data File : VX009461.D
 Acq On : 10 May 2019 04:59
 Operator : JC/SP
 Sample : K2782-16
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SS052MW352-190507

Quant Time: May 10 08:59:05 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 09 07:31:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	188913	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	302967	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	265290	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	112641	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	123909	48.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.10%	
35) Dibromofluoromethane	5.50	113	91935	48.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.54%	
50) Toluene-d8	8.71	98	376963	49.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.04%	
62) 4-Bromofluorobenzene	11.13	95	123890	44.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.68%	
Target Compounds						
19) Methyl tert-butyl Ether	3.20	73	7933	1.153	ug/l	92
27) cis-1,2-Dichloroethene	4.59	96	8288	3.612	ug/l	88
44) Trichloroethene	7.21	130	31957	15.171	ug/l	99
64) Tetrachloroethene	9.34	164	5706	3.136	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX050919\
Data File : VX009461.D
Acq On : 10 May 2019 04:59
Operator : JC/SP
Sample : K2782-16
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SS052MW352-190507

Quant Time: May 10 08:59:05 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260
QLast Update : Thu May 09 07:31:15 2019
Response via : Initial Calibration

