

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062819\
 Data File : VX010548.D
 Acq On : 28 Jun 2019 20:28
 Operator : JC/SP
 Sample : K3521-38
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PK-101

Quant Time: Jun 29 06:32:34 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	5.67	168	246651	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	382251	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	348726	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	162999	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	117740	50.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.82%	
35) Dibromofluoromethane	5.50	113	114449	48.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.88%	
50) Toluene-d8	8.71	98	452178	50.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.08%	
62) 4-Bromofluorobenzene	11.13	95	153392	47.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.96%	
Target Compounds						
16) Acetone	2.45	43	2228	1.922	ug/l #	87
19) Methyl tert-butyl Ether	3.20	73	3463	0.515	ug/l	94
65) Chlorobenzene	10.13	112	22234	3.140	ug/l	93
88) 1,4-Dichlorobenzene	12.10	146	4686	0.855	ug/l #	65
91) 1,2-Dichlorobenzene	12.39	146	1630	0.305	ug/l	83

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062819\
Data File : VX010548.D
Acq On : 28 Jun 2019 20:28
Operator : JC/SP
Sample : K3521-38
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PK-10I

Quant Time: Jun 29 06:32:34 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

