

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100819\
 Data File : VX012918.D
 Acq On : 09 Oct 2019 03:32
 Operator : JC/SP
 Sample : K5154-22
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-102I

Quant Time: Oct 09 07:42:09 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X100819W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 09 01:51:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	176857	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	302548	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	256319	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	96196	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	118553	53.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.26%	
35) Dibromofluoromethane	5.49	113	90105	51.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.78%	
50) Toluene-d8	8.71	98	356137	52.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.16%	
62) 4-Bromofluorobenzene	11.14	95	117540	44.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.34%	
Target Compounds						
16) Acetone	2.43	43	3140	2.695	ug/l	90
27) cis-1,2-Dichloroethene	4.59	96	2303	1.043	ug/l	96
44) Trichloroethene	7.20	130	11819	5.425	ug/l	93
64) Tetrachloroethene	9.33	164	20482	11.521	ug/l	96

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100819\
Data File : VX012918.D
Acq On : 09 Oct 2019 03:32
Operator : JC/SP
Sample : K5154-22
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-102I

Quant Time: Oct 09 07:42:09 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X100819W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 09 01:51:23 2019
Response via : Initial Calibration

