

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX041119\
 Data File : VX008854.D
 Acq On : 11 Apr 2019 16:07
 Operator : JC/SP
 Sample : K2349-17DL 4X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 M-19-190409DL

Quant Time: Apr 12 08:14:35 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X040319W.M
 Quant Title : SW846 8260
 QLast Update : Thu Apr 04 04:47:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	192220	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	311099	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	267533	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	113147	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	127744	45.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.12%	
35) Dibromofluoromethane	5.50	113	91124	45.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.62%	
50) Toluene-d8	8.71	98	387190	48.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.78%	
62) 4-Bromofluorobenzene	11.13	95	130170	46.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.70%	
Target Compounds						
9) 1,1,2-Trichlorotrifluoroet	2.38	101	3464	1.627	ug/l	97
27) cis-1,2-Dichloroethene	4.60	96	53953	19.523	ug/l	89
44) Trichloroethene	7.21	130	11089	4.700	ug/l	99
64) Tetrachloroethene	9.34	164	98281	49.583	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX041119\
Data File : VX008854.D
Acq On : 11 Apr 2019 16:07
Operator : JC/SP
Sample : K2349-17DL 4X
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
M-19-190409DL

Quant Time: Apr 12 08:14:35 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X040319W.M
Quant Title : SW846 8260
QLast Update : Thu Apr 04 04:47:41 2019
Response via : Initial Calibration

