

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX032019\
 Data File : VX008269.D
 Acq On : 21 Mar 2019 03:16
 Operator : JC/SP
 Sample : K1912-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 M3-T3-OZ(PRE)

Quant Time: Mar 22 08:01:36 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031919W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 19 13:15:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	320514	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	529208	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	451986	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	181434	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	229233	47.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.98%	
35) Dibromofluoromethane	5.50	113	170217	48.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.10%	
50) Toluene-d8	8.71	98	652195	48.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.12%	
62) 4-Bromofluorobenzene	11.13	95	209376	45.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.54%	
Target Compounds						
9) 1,1,2-Trichlorotrifluoroet	2.39	101	3820	0.927	ug/l #	84
16) Acetone	2.45	43	5847	2.136	ug/l	94
24) 1,1-Dichloroethane	3.70	63	4953	0.581	ug/l #	89
44) Trichloroethene	7.21	130	8849	1.985	ug/l	94

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX032019\
Data File : VX008269.D
Acq On : 21 Mar 2019 03:16
Operator : JC/SP
Sample : K1912-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
M3-T3-OZ(PRE)

Quant Time: Mar 22 08:01:36 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031919W.M
Quant Title : SW846 8260
QLast Update : Tue Mar 19 13:15:13 2019
Response via : Initial Calibration

