

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX041219\
 Data File : VX008909.D
 Acq On : 13 Apr 2019 00:52
 Operator : JC/SP
 Sample : K2311-15
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 0238-COMP-CS-3-ELTRT-DISS

Quant Time: Apr 13 06:35:22 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	5.67	168	184326	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	300522	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	251728	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	101606	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	124892	46.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.90%	
35) Dibromofluoromethane	5.50	113	90596	47.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.30%	
50) Toluene-d8	8.71	98	370804	48.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.94%	
62) 4-Bromofluorobenzene	11.13	95	117524	43.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.64%	
Target Compounds						
11) Tert butyl alcohol	3.04	59	3941	6.261	ug/l #	91
16) Acetone	2.44	43	240481	155.300	ug/l	97
18) Methyl Acetate	2.78	43	21468	5.716	ug/l	98
20) Methylene Chloride	2.85	84	30624	11.815	ug/l	94

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX041219\
Data File : VX008909.D
Acq On : 13 Apr 2019 00:52
Operator : JC/SP
Sample : K2311-15
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
0238-COMP-CS-3-ELTRT-DISS

Quant Time: Apr 13 06:35:22 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

