

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091118\
 Data File : VX004477.D
 Acq On : 11 Sep 2018 20:34
 Operator : JC/MD
 Sample : J4857-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY038PS034-180910

Quant Time: Sep 12 11:25:14 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 12 02:17:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	189492	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	303204	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	294951	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	161855	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	138088	56.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.04%	
35) Dibromofluoromethane	5.50	113	121553	43.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.62%	
50) Toluene-d8	8.72	98	442674	47.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.16%	
62) 4-Bromofluorobenzene	11.14	95	153929	46.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.70%	
Target Compounds						
16) Acetone	2.45	43	26461	22.19	ug/l	98
27) cis-1,2-Dichloroethene	4.59	96	7651	3.13	ug/l	94
44) Trichloroethene	7.21	130	955	0.34	ug/l	77
64) Tetrachloroethene	9.34	164	2066	0.56	ug/l	91

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091118\
Data File : VX004477.D
Acq On : 11 Sep 2018 20:34
Operator : JC/MD
Sample : J4857-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY038PS034-180910

Quant Time: Sep 12 11:25:14 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 12 02:17:23 2018
Response via : Initial Calibration

