

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX101018\
 Data File : VX005217.D
 Acq On : 10 Oct 2018 20:01
 Operator : JC/MD
 Sample : J5317-05 5X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE-122D1-20181004

Quant Time: Oct 11 07:37:11 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 03 04:00:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	175730	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	269880	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	273039	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	151590	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	132904	52.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.88%	
35) Dibromofluoromethane	5.51	113	113810	49.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.52%	
50) Toluene-d8	8.72	98	398051	52.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.70%	
62) 4-Bromofluorobenzene	11.14	95	142886	52.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.32%	
Target Compounds						
9) 1,1,2-Trichlorotrifluoroet	2.39	101	1418	0.74	ug/l	96
16) Acetone	2.45	43	2261	1.54	ug/l	98
27) cis-1,2-Dichloroethene	4.60	96	1376	0.59	ug/l	71
44) Trichloroethene	7.21	130	190107	79.15	ug/l	95

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX101018\
Data File : VX005217.D
Acq On : 10 Oct 2018 20:01
Operator : JC/MD
Sample : J5317-05 5X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RE-122D1-20181004

Quant Time: Oct 11 07:37:11 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 03 04:00:03 2018
Response via : Initial Calibration

