

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122418\
 Data File : VX006800.D
 Acq On : 24 Dec 2018 17:31
 Operator : JC/MD
 Sample : J6548-20
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 G-B

Quant Time: Dec 25 04:04:29 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 19 15:20:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	744999	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	1095075	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	1004633	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	506138	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	367962	45.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.66%	
35) Dibromofluoromethane	5.50	113	319816	46.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.36%	
50) Toluene-d8	8.71	98	1301950	49.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.74%	
62) 4-Bromofluorobenzene	11.13	95	471199	49.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.56%	
Target Compounds						
16) Acetone	2.45	43	29285	7.939	ug/l	93
18) Methyl Acetate	2.78	43	21120	1.947	ug/l	96
20) Methylene Chloride	2.86	84	36130	4.736	ug/l	96
43) Isopropyl Acetate	6.46	43	35600	2.479	ug/l	98
52) Toluene	8.79	92	18985	1.066	ug/l	96
68) m/p-Xylenes	10.36	106	16292	1.189	ug/l	92

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122418\
Data File : VX006800.D
Acq On : 24 Dec 2018 17:31
Operator : JC/MD
Sample : J6548-20
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
G-B

Quant Time: Dec 25 04:04:29 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260
QLast Update : Wed Dec 19 15:20:58 2018
Response via : Initial Calibration

