

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060321\
 Data File : VX022397.D
 Acq On : 04 Jun 2021 03:46
 Operator : JC/MD
 Sample : PB36780ZHE#12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB36780ZHE#12

Quant Time: Jun 04 05:41:39 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 19 14:23:10 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.562	168	60985	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	130042	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	117511	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	43136	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	53267	56.623	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 113.240%			
35) Dibromofluoromethane	5.397	113	38609	46.818	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 93.640%			
50) Toluene-d8	8.653	98	162039	51.009	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 102.020%			
62) 4-Bromofluorobenzene	11.085	95	56486	48.107	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 96.220%			
Target Compounds					Qvalue	
16) Acetone	2.392	43	11813	28.839	ug/l	91
18) Methyl Acetate	2.721	43	1406	1.291	ug/l #	74
20) Methylene Chloride	2.800	84	2379	2.616	ug/l	94
43) Isopropyl Acetate	6.348	43	11283	4.680	ug/l #	91

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060321\
Data File : VX022397.D
Acq On : 04 Jun 2021 03:46
Operator : JC/MD
Sample : PB36780ZHE#12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB36780ZHE#12

Quant Time: Jun 04 05:41:39 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260
QLast Update : Wed May 19 14:23:10 2021
Response via : Initial Calibration

