

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062521\
 Data File : VX022933.D
 Acq On : 25 Jun 2021 19:11
 Operator : JC/MD
 Sample : M2838-22
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DUP10-20210622

Quant Time: Jun 26 03:55:35 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 23 19:12:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	51297	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.769	114	109795	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	101609	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	44391	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.970	65	48492	52.090	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 104.180%	
35) Dibromofluoromethane	5.397	113	34240	49.109	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 98.220%	
50) Toluene-d8	8.653	98	135347	49.531	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 99.060%	
62) 4-Bromofluorobenzene	11.085	95	51364	50.366	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 100.740%	
Target Compounds						
						Qvalue
16) Acetone	2.386	43	708	1.715	ug/l	95
44) Trichloroethene	7.135	130	29418	36.889	ug/l	91
64) Tetrachloroethene	9.287	164	677	0.948	ug/l #	82

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062521\
Data File : VX022933.D
Acq On : 25 Jun 2021 19:11
Operator : JC/MD
Sample : M2838-22
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP10-20210622

Quant Time: Jun 26 03:55:35 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 23 19:12:26 2021
Response via : Initial Calibration

