

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060322\
 Data File : VX029198.D
 Acq On : 03 Jun 2022 18:22
 Operator : JC/MD
 Sample : N3164-13
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 37BR-20220531

Quant Time: Jun 06 02:36:32 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 01 04:45:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	303361	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	561868	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	545747	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	250674	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	199280	58.255	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	116.500%
35) Dibromofluoromethane	5.391	113	185437	49.688	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	99.380%
50) Toluene-d8	8.653	98	666578	48.835	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	97.660%
62) 4-Bromofluorobenzene	11.079	95	262348	49.722	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	99.440%
Target Compounds						
					Qvalue	
3) Chloromethane	1.288	50	642	0.233	ug/l #	84
16) Acetone	2.380	43	5611	4.165	ug/l	95
27) cis-1,2-Dichloroethene	4.489	96	2409	0.655	ug/l	91
37) Ethyl Acetate	4.721	43	146789	31.541	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060322\
Data File : VX029198.D
Acq On : 03 Jun 2022 18:22
Operator : JC/MD
Sample : N3164-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
37BR-20220531

Quant Time: Jun 06 02:36:32 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053122W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 01 04:45:42 2022
Response via : Initial Calibration

