

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
 Data File : VX033132.D
 Acq On : 30 Nov 2022 15:19
 Operator : JC/MD
 Sample : N5809-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 HOWE-DRUM-B-1129

Quant Time: Dec 01 00:20:48 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 22 13:58:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	140118	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	226770	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	212801	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	97155	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	76237	56.061	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	112.120%
35) Dibromofluoromethane	5.385	113	59385	46.451	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.900%
50) Toluene-d8	8.653	98	209929	54.782	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	109.560%
62) 4-Bromofluorobenzene	11.079	95	77390	46.083	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.160%
Target Compounds						
					Qvalue	
16) Acetone	2.398	43	14962	23.478	ug/l	98
20) Methylene Chloride	2.794	84	3812	1.619	ug/l	91
37) Ethyl Acetate	4.721	43	20559	11.528	ug/l	99
43) Isopropyl Acetate	6.342	43	88235	29.916	ug/l	100

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033132.D
Acq On : 30 Nov 2022 15:19
Operator : JC/MD
Sample : N5809-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
HOWE-DRUM-B-1129

Quant Time: Dec 01 00:20:48 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 22 13:58:46 2022
Response via : Initial Calibration

