

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060622\
 Data File : VX029272.D
 Acq On : 07 Jun 2022 05:31
 Operator : JC/MD
 Sample : N3166-24
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 19BR-20220601

Quant Time: Jun 07 10:07:38 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	5.556	168	303893	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	572274	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	567004	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	262368	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	214648	62.638	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	125.280%
35) Dibromofluoromethane	5.391	113	192994	50.772	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	101.540%
50) Toluene-d8	8.653	98	684619	49.244	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	98.480%
62) 4-Bromofluorobenzene	11.079	95	276240	51.403	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	102.800%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	7920	5.869	ug/l	92
27) cis-1,2-Dichloroethene	4.495	96	13688	3.714	ug/l	86
29) Tetrahydrofuran	5.026	42	2765	1.981	ug/l #	69
37) Ethyl Acetate	4.721	43	93764	19.781	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060622\
Data File : VX029272.D
Acq On : 07 Jun 2022 05:31
Operator : JC/MD
Sample : N3166-24
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
19BR-20220601

Quant Time: Jun 07 10:07:38 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

