

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032222\
 Data File : VN071475.D
 Acq On : 22 Mar 2022 18:33
 Operator : JC\MD
 Sample : N2031-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FIELD-BLANK

Quant Time: Mar 23 07:10:24 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 08 06:52:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	746585	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1281017	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1209798	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	429714	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	498838	48.752	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	97.500%
35) Dibromofluoromethane	8.016	113	374640	46.345	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	92.700%
50) Toluene-d8	10.439	98	1570442	48.251	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.500%
62) 4-Bromofluorobenzene	12.727	95	569556	51.678	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	103.360%
Target Compounds						
					Qvalue	
16) Acetone	4.298	43	9052	1.818	ug/l	99
18) Methyl Acetate	4.857	43	11475	1.320	ug/l	94
20) Methylene Chloride	5.098	84	4225	0.843	ug/l #	81

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032222\
Data File : VN071475.D
Acq On : 22 Mar 2022 18:33
Operator : JC\MD
Sample : N2031-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FIELD-BLANK

Quant Time: Mar 23 07:10:24 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
Quant Title : SW846 8260
QLast Update : Tue Mar 08 06:52:46 2022
Response via : Initial Calibration

