

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN079002.D
 Acq On : 22 Aug 2023 18:24
 Operator : JC\MD
 Sample : 04090-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-L

Quant Time: Aug 23 03:06:58 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	180229	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	350501	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	348461	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	109083	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	142846	50.934	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	101.860%
35) Dibromofluoromethane	8.171	113	119677	50.724	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.440%
50) Toluene-d8	10.571	98	445330	50.833	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	101.660%
62) 4-Bromofluorobenzene	12.853	95	135715	47.395	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.780%
Target Compounds						
						Qvalue
16) Acetone	4.442	43	25563	22.091	ug/l	91
20) Methylene Chloride	5.277	84	13268	5.380	ug/l #	78
43) Isopropyl Acetate	8.694	43	187816	26.760	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	40322	9.686	ug/l #	76

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
Data File : VN079002.D
Acq On : 22 Aug 2023 18:24
Operator : JC\MD
Sample : 04090-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-L

Quant Time: Aug 23 03:06:58 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
Quant Title : SW846 8260
QLast Update : Wed Aug 16 02:42:02 2023
Response via : Initial Calibration

