

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042023\
 Data File : VN077398.D
 Acq On : 20 Apr 2023 14:21
 Operator : JC\MD
 Sample : 02389-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SPB-2WC-23-04-17

Quant Time: Apr 21 02:39:47 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N040523W.M
 Quant Title : SW846 8260
 QLast Update : Thu Apr 06 08:27:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.231	168	540886	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.107	114	1004781	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.872	117	897822	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.795	152	318252	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.584	65	334408	45.236	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	90.480%
35) Dibromofluoromethane	8.172	113	303802	47.231	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.460%
50) Toluene-d8	10.572	98	1211294	46.778	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	93.560%
62) 4-Bromofluorobenzene	12.854	95	398564	46.280	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.560%
Target Compounds						
					Qvalue	
16) Acetone	4.449	43	42858	15.724	ug/l	98
20) Methylene Chloride	5.290	84	37632	4.889	ug/l	91
37) Ethyl Acetate	7.572	43	85700	10.068	ug/l	96
43) Isopropyl Acetate	8.695	43	397850	25.331	ug/l #	89

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042023\
Data File : VN077398.D
Acq On : 20 Apr 2023 14:21
Operator : JC\MD
Sample : 02389-06
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SPB-2WC-23-04-17

Quant Time: Apr 21 02:39:47 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N040523W.M
Quant Title : SW846 8260
QLast Update : Thu Apr 06 08:27:03 2023
Response via : Initial Calibration

