

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030922\
 Data File : VN071221.D
 Acq On : 10 Mar 2022 08:58
 Operator : JC\MD
 Sample : N1830-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 11977

Quant Time: Mar 11 00:30:07 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	8.087	168	662729	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1084009	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	974511	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.669	152	319534	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	430965	47.448	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	94.900%
35) Dibromofluoromethane	8.016	113	320723	46.886	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	93.780%
50) Toluene-d8	10.439	98	1338135	48.586	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	97.180%
62) 4-Bromofluorobenzene	12.727	95	455940	48.887	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	97.780%
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	69506	15.727	ug/l	96
18) Methyl Acetate	4.869	43	24393	2.506	ug/l	93
20) Methylene Chloride	5.098	84	27450	3.754	ug/l	91
43) Isopropyl Acetate	8.557	43	202590	8.982	ug/l #	69

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030922\
Data File : VN071221.D
Acq On : 10 Mar 2022 08:58
Operator : JC\MD
Sample : N1830-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
11977

Quant Time: Mar 11 00:30:07 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

