

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011222\
 Data File : VN070591.D
 Acq On : 12 Jan 2022 17:17
 Operator : JC/MD
 Sample : N1097-13
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-3S

Quant Time: Jan 13 04:27:36 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
 Quant Title : SW846 8260
 QLast Update : Mon Dec 27 18:47:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.083	168	841547	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.965	114	1485717	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1332173	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	464351	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	656908	54.291	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	108.580%
35) Dibromofluoromethane	8.021	113	462573	51.458	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	102.920%
50) Toluene-d8	10.440	98	1887194	51.391	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	102.780%
62) 4-Bromofluorobenzene	12.731	95	693425	52.281	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	104.560%
Target Compounds						
						Qvalue
16) Acetone	4.291	43	58943	7.504	ug/l	94
18) Methyl Acetate	4.859	43	36628	1.612	ug/l	93
20) Methylene Chloride	5.098	84	32482	2.989	ug/l	93
43) Isopropyl Acetate	8.560	43	162534	5.348	ug/l #	82

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011222\
Data File : VN070591.D
Acq On : 12 Jan 2022 17:17
Operator : JC/MD
Sample : N1097-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-3S

Quant Time: Jan 13 04:27:36 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260
QLast Update : Mon Dec 27 18:47:12 2021
Response via : Initial Calibration

