

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102121\
 Data File : VN069118.D
 Acq On : 22 Oct 2021 4:37
 Operator : JC/MD
 Sample : M4285-13
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-GB-(1)

Quant Time: Oct 22 07:21:43 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	336379	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	583957	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	588337	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	217429	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	216921	48.911	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	97.820%
35) Dibromofluoromethane	8.024	113	170850	50.287	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	100.580%
50) Toluene-d8	10.443	98	735466	54.102	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	108.200%
62) 4-Bromofluorobenzene	12.733	95	265672	51.799	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	103.600%
Target Compounds						
					Qvalue	
16) Acetone	4.301	43	12316	7.441	ug/l	98
18) Methyl Acetate	4.867	43	8776	1.398	ug/l	98
20) Methylene Chloride	5.103	84	12020	3.155	ug/l	92
43) Isopropyl Acetate	8.566	43	58439	6.429	ug/l #	79

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102121\
Data File : VN069118.D
Acq On : 22 Oct 2021 4:37
Operator : JC/MD
Sample : M4285-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MH-GB-(1)

Quant Time: Oct 22 07:21:43 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 13 05:44:40 2021
Response via : Initial Calibration

