

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102121\
 Data File : VN069115.D
 Acq On : 22 Oct 2021 3:22
 Operator : JC/MD
 Sample : M4285-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 41 Sample Multiplier: 1

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

Quant Time: Oct 22 07:21:13 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	361676	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	629669	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	623257	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	240911	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	232224	48.699	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	97.400%
35) Dibromofluoromethane	8.026	113	180916	49.384	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	98.760%
50) Toluene-d8	10.446	98	785628	53.597	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	107.200%
62) 4-Bromofluorobenzene	12.733	95	286845	51.867	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	103.740%
Target Compounds						
					Qvalue	
16) Acetone	4.296	43	11414	6.414	ug/l	97
18) Methyl Acetate	4.862	43	8410	1.246	ug/l #	75
20) Methylene Chloride	5.098	84	11746	2.868	ug/l	93
43) Isopropyl Acetate	8.568	43	58671	5.986	ug/l #	80

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102121\
Data File : VN069115.D
Acq On : 22 Oct 2021 3:22
Operator : JC/MD
Sample : M4285-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 41 Sample Multiplier: 1

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

Quant Time: Oct 22 07:21:13 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

