

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122922\
 Data File : VN076046.D
 Acq On : 29 Dec 2022 14:06
 Operator : JC\MD
 Sample : N6228-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-2

Quant Time: Dec 30 05:32:50 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 14 13:03:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	364074	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	565397	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	511182	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	205914	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	178295	45.015	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	90.040%
35) Dibromofluoromethane	8.177	113	168693	48.178	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.360%
50) Toluene-d8	10.571	98	661055	50.495	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	101.000%
62) 4-Bromofluorobenzene	12.847	95	206917	47.720	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.440%
Target Compounds						
					Qvalue	
16) Acetone	4.453	43	42631	31.223	ug/l	99
20) Methylene Chloride	5.294	84	13134	2.994	ug/l #	92
37) Ethyl Acetate	7.577	43	31831	8.520	ug/l	99
43) Isopropyl Acetate	8.694	43	188027	30.509	ug/l #	83

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122922\
Data File : VN076046.D
Acq On : 29 Dec 2022 14:06
Operator : JC\MD
Sample : N6228-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-2

Quant Time: Dec 30 05:32:50 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260
QLast Update : Wed Dec 14 13:03:41 2022
Response via : Initial Calibration

