

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081522\
 Data File : VN073921.D
 Acq On : 16 Aug 2022 02:31
 Operator : JC\MD
 Sample : N4188-11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 S15

Quant Time: Aug 16 05:14:23 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081522W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 16 00:45:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	640645	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	929135	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	906406	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	492096	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	298311	47.807	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	95.620%
35) Dibromofluoromethane	7.951	113	290276	52.827	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	105.660%
50) Toluene-d8	10.380	98	989851	46.513	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	93.020%
62) 4-Bromofluorobenzene	12.674	95	416735	59.005	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	118.000%
Target Compounds						
					Qvalue	
16) Acetone	4.222	43	90455	30.726	ug/l	98
20) Methylene Chloride	5.022	84	30910	4.976	ug/l	88
37) Ethyl Acetate	7.339	43	69140	7.478	ug/l	98
43) Isopropyl Acetate	8.492	43	337633	25.633	ug/l #	90

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081522\
Data File : VN073921.D
Acq On : 16 Aug 2022 02:31
Operator : JC\MD
Sample : N4188-11
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
S15

Quant Time: Aug 16 05:14:23 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081522W.M
Quant Title : SW846 8260
QLast Update : Tue Aug 16 00:45:21 2022
Response via : Initial Calibration

