

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122022\
 Data File : VN075896.D
 Acq On : 20 Dec 2022 14:11
 Operator : JC\MD
 Sample : N6129-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 C3

Quant Time: Dec 21 01:28:10 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	8.235	168	417598	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	656067	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	572281	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	245810	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	207582	45.692	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	91.380%
35) Dibromofluoromethane	8.177	113	191344	47.095	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.200%
50) Toluene-d8	10.571	98	761047	50.099	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	100.200%
62) 4-Bromofluorobenzene	12.853	95	247135	49.119	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.240%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	60382	38.556	ug/l	100
20) Methylene Chloride	5.294	84	21239	4.487	ug/l	95
37) Ethyl Acetate	7.571	43	37311	8.606	ug/l	99
43) Isopropyl Acetate	8.694	43	251861	35.219	ug/l #	80
51) 4-Methyl-2-Pentanone	10.447	43	24547	5.676	ug/l #	79

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

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