

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122022\
 Data File : VN075897.D
 Acq On : 20 Dec 2022 14:35
 Operator : JC\MD
 Sample : N6129-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 D8

Quant Time: Dec 21 01:28:21 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	420091	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	658771	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	581134	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	248968	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	211573	46.294	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	92.580%
35) Dibromofluoromethane	8.171	113	193642	47.465	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.940%
50) Toluene-d8	10.571	98	761562	49.927	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.860%
62) 4-Bromofluorobenzene	12.853	95	246020	48.697	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.400%
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	58303	37.008	ug/l	96
20) Methylene Chloride	5.294	84	20440	4.264	ug/l #	95
37) Ethyl Acetate	7.571	43	36154	8.305	ug/l	96
43) Isopropyl Acetate	8.694	43	372549	51.881	ug/l #	71
51) 4-Methyl-2-Pentanone	10.453	43	25182	5.799	ug/l #	79

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

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