

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122822\
 Data File : VN076022.D
 Acq On : 29 Dec 2022 00:55
 Operator : JC\MD
 Sample : PB149923ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB149923ZHE#03

Quant Time: Dec 29 07:49:02 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	387000	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	608242	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	549743	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	232835	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	195886	46.527	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	93.060%	
35) Dibromofluoromethane	8.177	113	177540	47.133	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	94.260%	
50) Toluene-d8	10.571	98	719485	51.087	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	102.180%	
62) 4-Bromofluorobenzene	12.853	95	231444	49.617	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	99.240%	
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	68693	47.331	ug/l	98
20) Methylene Chloride	5.295	84	18965	4.299	ug/l #	86
25) 2-Butanone	7.489	43	25120	12.361	ug/l	91
37) Ethyl Acetate	7.571	43	35663	8.873	ug/l	98
43) Isopropyl Acetate	8.694	43	335349	50.580	ug/l #	73
51) 4-Methyl-2-Pentanone	10.447	43	22149	5.524	ug/l #	81

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

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