

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090223\
 Data File : VN079141.D
 Acq On : 02 Sep 2023 14:23
 Operator : JC\MD
 Sample : PB155301ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155301ZHE#03

Quant Time: Sep 04 01:03:40 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	216781	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	413338	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	405596	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	117634	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	164042	48.630	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.260%
35) Dibromofluoromethane	8.171	113	140813	50.609	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.220%
50) Toluene-d8	10.571	98	537268	52.004	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	104.000%
62) 4-Bromofluorobenzene	12.853	95	159863	47.341	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.680%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	29006	20.840	ug/l	92
20) Methylene Chloride	5.289	84	28913	9.748	ug/l	89
37) Ethyl Acetate	7.571	43	4879	1.042	ug/l #	79
43) Isopropyl Acetate	8.694	43	184077	22.240	ug/l #	83
51) 4-Methyl-2-Pentanone	10.453	43	30805	6.275	ug/l #	83

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

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