

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090223\
 Data File : VN079138.D
 Acq On : 02 Sep 2023 13:11
 Operator : JC\MD
 Sample : PB155301TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155301TB

Quant Time: Sep 04 01:02:17 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	277068	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	509411	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	483766	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	147584	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	157654	36.567	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	73.140%#
35) Dibromofluoromethane	8.171	113	176083	51.350	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.700%
50) Toluene-d8	10.571	98	633824	49.780	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.560%
62) 4-Bromofluorobenzene	12.853	95	140908	33.858	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	67.720%#
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	29132	16.376	ug/l	91
20) Methylene Chloride	5.283	84	21285	5.615	ug/l #	88
43) Isopropyl Acetate	8.694	43	124571	12.212	ug/l #	83
51) 4-Methyl-2-Pentanone	10.447	43	33204	5.488	ug/l #	79

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

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