

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090223\
 Data File : VN079144.D
 Acq On : 02 Sep 2023 15:35
 Operator : JC\MD
 Sample : PB155301ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155301ZHE#06

Quant Time: Sep 04 01:04:57 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	207740	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	290153	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	401625	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	120780	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	161459	49.947	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.900%
35) Dibromofluoromethane	8.171	113	135711	69.483	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	138.960%#
50) Toluene-d8	10.571	98	526928	72.657	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	145.320%#
62) 4-Bromofluorobenzene	12.853	95	161070	67.949	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	135.900%#
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	29662	22.238	ug/l	97
20) Methylene Chloride	5.271	84	21208	7.461	ug/l #	91
37) Ethyl Acetate	7.577	43	5052	1.537	ug/l #	87
43) Isopropyl Acetate	8.694	43	225746	38.853	ug/l #	77
51) 4-Methyl-2-Pentanone	10.447	43	31637	9.181	ug/l #	85

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

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