

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082923\
 Data File : VN079092.D
 Acq On : 29 Aug 2023 14:29
 Operator : JC\MD
 Sample : PB155153ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155153ZHE#01

Quant Time: Aug 29 23:29:41 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	144307	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	264162	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	269103	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	76481	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	110967	49.417	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	98.840%
35) Dibromofluoromethane	8.171	113	94341	53.054	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	106.100%
50) Toluene-d8	10.571	98	353549	53.547	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	107.100%
62) 4-Bromofluorobenzene	12.853	95	101195	46.890	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.780%
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	32891	35.499	ug/l #	87
20) Methylene Chloride	5.289	84	20593	10.429	ug/l #	84
25) 2-Butanone	7.494	43	18246	12.283	ug/l #	79
43) Isopropyl Acetate	8.694	43	154692	29.244	ug/l #	79
51) 4-Methyl-2-Pentanone	10.453	43	62268	19.847	ug/l	87

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

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