

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090223\
 Data File : VN079145.D
 Acq On : 02 Sep 2023 15:59
 Operator : JC\MD
 Sample : PB155301ZHE#07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155301ZHE#07

Quant Time: Sep 04 01:05:23 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	8.229	168	209207	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	408476	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	401645	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	113697	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	161391	49.576	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.160%
35) Dibromofluoromethane	8.171	113	138040	50.203	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.400%
50) Toluene-d8	10.570	98	524799	51.402	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	102.800%
62) 4-Bromofluorobenzene	12.853	95	158287	47.432	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.860%
Target Compounds						
					Qvalue	
16) Acetone	4.441	43	28103	20.922	ug/l	95
20) Methylene Chloride	5.288	84	21095	7.369	ug/l #	82
37) Ethyl Acetate	7.577	43	4654	1.006	ug/l #	81
43) Isopropyl Acetate	8.694	43	229403	28.046	ug/l #	77
51) 4-Methyl-2-Pentanone	10.447	43	32020	6.600	ug/l #	79

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

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