

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090223\
 Data File : VN079139.D
 Acq On : 02 Sep 2023 13:35
 Operator : JC\MD
 Sample : PB155301ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155301ZHE#01

Quant Time: Sep 04 01:02:45 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	213592	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	407579	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	306753	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	120833	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	161153	48.486	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	96.980%
35) Dibromofluoromethane	8.171	113	139758	50.939	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.880%
50) Toluene-d8	10.571	98	534825	52.499	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	105.000%
62) 4-Bromofluorobenzene	12.853	95	163641	49.144	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.280%
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	27430	20.001	ug/l	99
20) Methylene Chloride	5.283	84	29871	10.221	ug/l	97
43) Isopropyl Acetate	8.694	43	138261	16.940	ug/l #	82
51) 4-Methyl-2-Pentanone	10.447	43	32099	6.631	ug/l #	79

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

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