

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082323\
 Data File : VN079012.D
 Acq On : 23 Aug 2023 13:35
 Operator : JC\MD
 Sample : PB155023TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155023TB

Quant Time: Aug 24 01:09:57 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	182641	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	337073	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	334670	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	102069	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	140525	49.445	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	98.880%
35) Dibromofluoromethane	8.171	113	119510	52.671	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	105.340%
50) Toluene-d8	10.571	98	438693	52.070	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	104.140%
62) 4-Bromofluorobenzene	12.853	95	135258	49.117	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.240%
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	30856	26.313	ug/l	96
20) Methylene Chloride	5.294	84	10658	4.265	ug/l #	89
25) 2-Butanone	7.488	43	17580	9.351	ug/l #	70
43) Isopropyl Acetate	8.694	43	165535	24.525	ug/l #	77
51) 4-Methyl-2-Pentanone	10.447	43	68862	17.201	ug/l #	77

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

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