

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100423\
 Data File : VN079403.D
 Acq On : 04 Oct 2023 15:11
 Operator : JC\MD
 Sample : PB156087ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB156087ZHE#01

Quant Time: Oct 05 02:21:43 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091123W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 12 05:05:24 2023
 Response via : Initial Calibration

Compound	R.T. QIon		Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.230	168	123999	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	244874	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	243420	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	57255	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	108099	64.580	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 129.160%#		
35) Dibromofluoromethane	8.171	113	89842	55.157	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 110.320%		
50) Toluene-d8	10.571	98	331474	55.422	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 110.840%		
62) 4-Bromofluorobenzene	12.853	95	84926	44.515	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 89.040%		
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	22537	31.702	ug/l	93
20) Methylene Chloride	5.283	84	16966	9.807	ug/l	87
43) Isopropyl Acetate	8.694	43	133313	32.626	ug/l #	80
51) 4-Methyl-2-Pentanone	10.447	43	13351	5.629	ug/l #	66

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

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