

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091523\
 Data File : VN079275.D
 Acq On : 15 Sep 2023 16:30
 Operator : JC\MD
 Sample : PB155573ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155573ZHE#06

Quant Time: Sep 16 01:45:07 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	8.229	168	144995	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	286741	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	287116	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	82145	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	121153	61.898	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	123.800%
35) Dibromofluoromethane	8.171	113	99800	52.324	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	104.640%
50) Toluene-d8	10.570	98	382668	54.640	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	109.280%
62) 4-Bromofluorobenzene	12.853	95	112268	50.255	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.500%
Target Compounds						
						Qvalue
16) Acetone	4.436	43	19381	23.315	ug/l	90
20) Methylene Chloride	5.277	84	12045	5.954	ug/l #	83
25) 2-Butanone	7.494	43	11351	9.411	ug/l #	69
43) Isopropyl Acetate	8.694	43	182044	38.047	ug/l #	76
51) 4-Methyl-2-Pentanone	10.447	43	31844	11.465	ug/l #	81

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

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