

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082436.D
 Acq On : 03 Jun 2024 16:00
 Operator : JC\MD
 Sample : PB161254ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161254ZHE#03

Quant Time: Jun 04 00:31:01 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 16 05:21:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.229	168	156740	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	289546	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	284736	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	121898	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	134416	54.310	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 108.620%	
35) Dibromofluoromethane	8.171	113	94993	53.332	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 106.660%	
50) Toluene-d8	10.570	98	376435	55.125	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 110.240%	
62) 4-Bromofluorobenzene	12.853	95	130876	53.696	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	= 107.400%	
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	85856	84.564	ug/l	92
20) Methylene Chloride	5.277	84	12788	6.243	ug/l	88
25) 2-Butanone	7.488	43	14476	9.404	ug/l #	73
43) Isopropyl Acetate	8.694	43	334434	56.456	ug/l #	70
51) 4-Methyl-2-Pentanone	10.447	43	20108	5.974	ug/l #	73

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

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