

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
 Data File : VN082638.D
 Acq On : 19 Jun 2024 13:57
 Operator : JC\MD
 Sample : PB161542ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161542ZHE#04

Quant Time: Jun 20 03:00:41 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061724W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 18 04:52:44 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	134349	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	253432	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	263099	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	100101	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	119573	55.500	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	111.000%
35) Dibromofluoromethane	8.171	113	88383	51.408	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.820%
50) Toluene-d8	10.570	98	343875	52.891	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	105.780%
62) 4-Bromofluorobenzene	12.853	95	110864	49.459	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	=	98.920%
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	57824	59.014	ug/l	92
20) Methylene Chloride	5.283	84	11642	6.758	ug/l	92
25) 2-Butanone	7.488	43	17645	12.670	ug/l	97
43) Isopropyl Acetate	8.694	43	251450	41.098	ug/l #	83
51) 4-Methyl-2-Pentanone	10.447	43	16553	5.275	ug/l #	82

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

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