

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
 Data File : VN082635.D
 Acq On : 19 Jun 2024 12:44
 Operator : JC\MD
 Sample : PB161542ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161542ZHE#01

Quant Time: Jun 20 02:59:10 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	8.229	168	138256	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	258732	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	266590	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	108848	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	123722	55.803	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	111.600%	
35) Dibromofluoromethane	8.171	113	91657	52.220	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	104.440%	
50) Toluene-d8	10.570	98	349242	52.616	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	105.240%	
62) 4-Bromofluorobenzene	12.853	95	113649	49.663	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	=	99.320%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	60409	59.910	ug/l	98
20) Methylene Chloride	5.271	84	11465	6.467	ug/l	98
25) 2-Butanone	7.494	43	18088	12.621	ug/l	99
37) Ethyl Acetate	7.577	43	4870	1.457	ug/l #	88
43) Isopropyl Acetate	8.694	43	224540	35.948	ug/l #	85
51) 4-Methyl-2-Pentanone	10.453	43	16829	5.253	ug/l #	78

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

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