

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061024\
 Data File : VN082511.D
 Acq On : 10 Jun 2024 13:06
 Operator : JC\MD
 Sample : PB161359TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161359TB

Manual Integrations
 APPROVED

Reviewed By :Mahesh Dadoda 06/11/2024
 Supervised By :Semsettin Yesilyurt 06/11/2024

Quant Time: Jun 11 01:06:00 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	8.230	168	176405	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	321122	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	322242	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	131071	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	137028	49.194	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	98.380%		
35) Dibromofluoromethane	8.165	113	100162	50.705	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	101.400%		
50) Toluene-d8	10.565	98	389759	51.463	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	102.920%		
62) 4-Bromofluorobenzene	12.853	95	130927	48.435	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery =	96.860%		
Target Compounds					Qvalue	
16) Acetone	4.430	43	96988	84.880	ug/l #	88
20) Methylene Chloride	5.283	84	6339	2.750	ug/l #	77
25) 2-Butanone	7.488	43	24712	14.264	ug/l #	87
37) Ethyl Acetate	7.571	43	5333	1.438	ug/l #	82
43) Isopropyl Acetate	8.688	43	227050	34.560	ug/l #	76
51) 4-Methyl-2-Pentanone	10.447	43	23873m	6.396	ug/l	

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

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