

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010423\
 Data File : VN076086.D
 Acq On : 04 Jan 2023 17:06
 Operator : JC\MD
 Sample : PB150013ZHE#07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB150013ZHE#07

Manual Integrations
 APPROVED

Reviewed By : John Carlone 01/05/2023
 Supervised By : Mahesh Dadoda 01/05/2023

Quant Time: Jan 05 01:25:44 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 14 13:03:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.236	168	328722	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	534907	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	478851	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	200252	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.589	65	173634	48.553	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	97.100%		
35) Dibromofluoromethane	8.177	113	158103	47.728	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.460%		
50) Toluene-d8	10.571	98	625609	50.512	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	101.020%		
62) 4-Bromofluorobenzene	12.853	95	197330	48.104	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	96.200%		
Target Compounds					Qvalue	
16) Acetone	4.442	43	41621	33.762	ug/l	95
20) Methylene Chloride	5.289	84	13755	3.576	ug/l	98
25) 2-Butanone	7.500	43	8623	4.996	ug/l	89
37) Ethyl Acetate	7.577	43	30843	8.726	ug/l	99
43) Isopropyl Acetate	8.700	43	262480m	45.017	ug/l	

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

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