

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010423\
 Data File : VN076083.D
 Acq On : 04 Jan 2023 15:53
 Operator : JC\MD
 Sample : PB150013ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB150013ZHE#04

Quant Time: Jan 05 01:24:42 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

Manual Integrations
 APPROVED

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

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	325707	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	537200	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	481352	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	191797	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	175783	49.609	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	99.220%		
35) Dibromofluoromethane	8.177	113	157074	47.215	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.420%		
50) Toluene-d8	10.571	98	628235	50.507	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	101.020%		
62) 4-Bromofluorobenzene	12.853	95	198127	48.092	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	96.180%		
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	41596	34.054	ug/l	94
20) Methylene Chloride	5.283	84	11239	2.836	ug/l #	78
25) 2-Butanone	7.494	43	9403	5.498	ug/l #	87
37) Ethyl Acetate	7.571	43	30383	8.559	ug/l	98
43) Isopropyl Acetate	8.694	43	256652m	43.830	ug/l	

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

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