

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122322\
 Data File : VN075966.D
 Acq On : 23 Dec 2022 15:09
 Operator : JC\MD
 Sample : PB149839TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB149839TB

Manual Integrations
 APPROVED

Reviewed By :Krupa Patel 12/27/2022
 Supervised By :Mahesh Dadoda 12/27/2022

Quant Time: Dec 26 04:33:54 2022
 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	340063	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	552511	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	482039	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	196967	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	181250	48.992	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	97.980%		
35) Dibromofluoromethane	8.177	113	161399	47.170	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.340%		
50) Toluene-d8	10.571	98	646796	50.558	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	101.120%		
62) 4-Bromofluorobenzene	12.853	95	199539	47.092	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	94.180%		
Target Compounds						Qvalue
16) Acetone	4.442	43	43614	34.199	ug/l	93
20) Methylene Chloride	5.295	84	16873	4.361	ug/l	93
25) 2-Butanone	7.500	43	11333	6.347	ug/l #	86
37) Ethyl Acetate	7.571	43	32937	9.021	ug/l	99
43) Isopropyl Acetate	8.694	43	254822m	42.311	ug/l	

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

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