

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112322\
 Data File : VN075398.D
 Acq On : 23 Nov 2022 15:59
 Operator : JC\MD
 Sample : PB149245ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB149245ZHE#01

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2022
 Supervised By : Mahesh Dadoda 11/28/2022

Quant Time: Nov 24 06:52:29 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 23 00:18:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.229	168	162307	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	277919	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	258719	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	118331	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	44755	51.369	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 102.740%			
35) Dibromofluoromethane	8.171	113	54686	50.090	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.180%			
50) Toluene-d8	10.570	98	130040	52.885	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 105.760%			
62) 4-Bromofluorobenzene	12.853	95	93130	51.312	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 102.620%			
Target Compounds						
8) Diethyl Ether	3.965	74	4296m	4.408	ug/l	
16) Acetone	4.441	43	81812	100.673	ug/l	89
20) Methylene Chloride	5.294	84	21881	12.314	ug/l	95
37) Ethyl Acetate	7.565	43	23220	10.831	ug/l	97
43) Isopropyl Acetate	8.694	43	119656	34.773	ug/l #	90

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

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