

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112222\
 Data File : VN075367.D
 Acq On : 22 Nov 2022 14:56
 Operator : JC\MD
 Sample : PB149227TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB149227TB

Quant Time: Nov 23 07:09:36 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	141354	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	244746	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	225420	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	99092	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	40431	53.285	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 106.580%	
35) Dibromofluoromethane	8.171	113	49334	51.312	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.620%	
50) Toluene-d8	10.565	98	112822	52.101	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 104.200%	
62) 4-Bromofluorobenzene	12.847	95	80304	50.243	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 100.480%	
Target Compounds						
					Qvalue	
8) Diethyl Ether	3.965	74	4290	5.055	ug/l	91
16) Acetone	4.441	43	76325	107.842	ug/l	96
20) Methylene Chloride	5.288	84	21144	13.807	ug/l #	91
37) Ethyl Acetate	7.571	43	22018m	11.663	ug/l	
43) Isopropyl Acetate	8.694	43	98856	32.622	ug/l #	91
51) 4-Methyl-2-Pentanone	10.447	43	8437	4.371	ug/l #	69

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

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