

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122022\
 Data File : VN075895.D
 Acq On : 20 Dec 2022 13:47
 Operator : JC\MD
 Sample : PB149742TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB149742TB

Quant Time: Dec 21 01:27:59 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	8.229	168	413870	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	642533	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	570035	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	251349	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	204450	45.408	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	90.820%
35) Dibromofluoromethane	8.171	113	187841	47.207	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.420%
50) Toluene-d8	10.570	98	740342	49.763	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.520%
62) 4-Bromofluorobenzene	12.853	95	243073	49.329	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.660%
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	57637	37.135	ug/l	98
20) Methylene Chloride	5.300	84	19078	4.006	ug/l	99
37) Ethyl Acetate	7.571	43	38690	9.112	ug/l	96
43) Isopropyl Acetate	8.694	43	225320	32.171	ug/l #	84
51) 4-Methyl-2-Pentanone	10.447	43	23851	5.631	ug/l #	81

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

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