

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072523\
 Data File : VN078645.D
 Acq On : 25 Jul 2023 14:23
 Operator : JC\MD
 Sample : PB154393TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154393TB

Quant Time: Jul 26 04:46:24 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.236	168	429455	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	752667	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	724223	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	235788	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	263878	45.834	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	91.660%	
35) Dibromofluoromethane	8.177	113	267569	52.369	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	104.740%	
50) Toluene-d8	10.571	98	931986	48.286	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	96.580%	
62) 4-Bromofluorobenzene	12.853	95	301765	50.757	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	101.520%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	82417	45.956	ug/l	99
20) Methylene Chloride	5.289	84	23139	3.357	ug/l	94
25) 2-Butanone	7.500	43	57653	20.422	ug/l #	79
37) Ethyl Acetate	7.571	43	18160	2.953	ug/l	97
43) Isopropyl Acetate	8.700	43	307455	30.365	ug/l #	86
51) 4-Methyl-2-Pentanone	10.453	43	193493	32.128	ug/l	92

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

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