

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN073123\
 Data File : VN078703.D
 Acq On : 31 Jul 2023 14:23
 Operator : JC\MD
 Sample : PB154517TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154517TB

Quant Time: Aug 01 01:17:30 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	723297	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	1281852	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	1208505	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	429018	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	469597	48.429	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	96.860%
35) Dibromofluoromethane	8.177	113	444452	51.077	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.160%
50) Toluene-d8	10.571	98	1638378	49.841	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.680%
62) 4-Bromofluorobenzene	12.853	95	556724	54.984	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	109.960%
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	176040	58.282	ug/l	94
20) Methylene Chloride	5.289	84	37977	3.240	ug/l	93
25) 2-Butanone	7.494	43	99762	20.982	ug/l #	81
37) Ethyl Acetate	7.571	43	36851	3.518	ug/l	98
43) Isopropyl Acetate	8.694	43	566882	32.874	ug/l #	86
51) 4-Methyl-2-Pentanone	10.453	43	422477	41.190	ug/l	92

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

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