

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080823\
 Data File : VN078796.D
 Acq On : 08 Aug 2023 12:47
 Operator : JC\MD
 Sample : PB154681TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154681TB

Quant Time: Aug 09 02:48:03 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080723W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 08 05:03:52 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	173511	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	315238	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	292373	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	77835	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	128468	49.388	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	98.780%
35) Dibromofluoromethane	8.177	113	102229	48.021	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.040%
50) Toluene-d8	10.571	98	394086	50.304	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	100.600%
62) 4-Bromofluorobenzene	12.853	95	111249	47.516	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.040%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	30082	30.057	ug/l	99
20) Methylene Chloride	5.271	84	12438	5.018	ug/l #	88
25) 2-Butanone	7.500	43	14979	8.838	ug/l #	66
43) Isopropyl Acetate	8.694	43	129600	21.890	ug/l #	81
51) 4-Methyl-2-Pentanone	10.447	43	56078	15.464	ug/l #	81

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

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