

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062223\
 Data File : VN078355.D
 Acq On : 22 Jun 2023 17:16
 Operator : JC\MD
 Sample : PB153671TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB153671TB

Quant Time: Jun 23 08:03:25 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060623W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 07 05:05:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	497645	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	911275	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	854153	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	281002	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	338409	51.003	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 102.000%	
35) Dibromofluoromethane	8.177	113	311580	54.076	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 108.160%	
50) Toluene-d8	10.571	98	1109745	49.960	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 99.920%	
62) 4-Bromofluorobenzene	12.853	95	366227	53.035	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 106.080%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	188662	98.784	ug/l	98
20) Methylene Chloride	5.295	84	22815	3.615	ug/l #	91
25) 2-Butanone	7.494	43	64927	22.107	ug/l	90
37) Ethyl Acetate	7.577	43	71416	11.061	ug/l	98
43) Isopropyl Acetate	8.700	43	335417	27.165	ug/l #	85
51) 4-Methyl-2-Pentanone	10.447	43	445199	68.470	ug/l	92

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

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