

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120924\
 Data File : VN085152.D
 Acq On : 09 Dec 2024 14:59
 Operator : JC\MD
 Sample : PB165391TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB165391TB

Quant Time: Dec 10 05:02:28 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112224W.M
 Quant Title : SW846 8260
 QLast Update : Sat Nov 23 02:54:10 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	135385	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	250920	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	218767	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	78597	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	103828	53.840	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 107.680%			
35) Dibromofluoromethane	8.165	113	84687	50.048	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.100%			
50) Toluene-d8	10.565	98	299518	49.392	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 98.780%			
62) 4-Bromofluorobenzene	12.847	95	98715	43.529	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 87.060%			
Target Compounds						
16) Acetone	4.436	43	37678	67.085	ug/l	95
20) Methylene Chloride	5.271	84	2457	1.434	ug/l #	80
43) Isopropyl Acetate	8.688	43	208335	48.136	ug/l #	78
51) 4-Methyl-2-Pentanone	10.441	43	10022	5.070	ug/l #	73

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

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