

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN092724\
 Data File : VN084201.D
 Acq On : 27 Sep 2024 14:18
 Operator : JC\MD
 Sample : PB163689TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB163689TB

Quant Time: Sep 28 00:59:59 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 11 15:18:56 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	150157	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	269344	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	229207	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	84285	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.576	65	112872	49.852	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	99.700%		
35) Dibromofluoromethane	8.165	113	85958	49.851	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	99.700%		
50) Toluene-d8	10.565	98	326357	54.749	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	109.500%		
62) 4-Bromofluorobenzene	12.847	95	102101	41.515	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	83.040%		
Target Compounds						
16) Acetone	4.430	43	57977	63.876	ug/l	96
20) Methylene Chloride	5.288	84	8299	4.468	ug/l #	79
25) 2-Butanone	7.482	43	8859	7.426	ug/l #	87
43) Isopropyl Acetate	8.688	43	147556	34.564	ug/l #	86
51) 4-Methyl-2-Pentanone	10.441	43	13242	5.698	ug/l #	78

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

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