

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082724\
 Data File : VN083510.D
 Acq On : 27 Aug 2024 15:29
 Operator : JC\MD
 Sample : PB162992ZHE#08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB162992ZHE#08

Manual Integrations
 APPROVED

Reviewed By : John Carlone 08/28/2024
 Supervised By : Mahesh Dadoda 08/28/2024

Quant Time: Aug 28 01:47:31 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080724W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 08 06:30:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	116057	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	245294	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	275564	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	135306	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	112622	68.176	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 136.360%#			
35) Dibromofluoromethane	8.165	113	88244	57.635	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 115.280%			
50) Toluene-d8	10.565	98	327773	57.391	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 114.780%#			
62) 4-Bromofluorobenzene	12.847	95	155235	69.719	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 139.440%#			
Target Compounds						Qvalue
16) Acetone	4.436	43	22278	34.465	ug/l	96
18) Methyl Acetate	5.018	43	11746	6.099	ug/l	96
43) Isopropyl Acetate	8.688	43	359832m	83.002	ug/l	

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

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