

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061024\
 Data File : VN082513.D
 Acq On : 10 Jun 2024 13:54
 Operator : JC\MD
 Sample : P2740-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WCS-TPB

Manual Integrations
 APPROVED

Reviewed By :Mahesh Dadoda 06/11/2024
 Supervised By :Semsettin Yesilyurt 06/11/2024

Quant Time: Jun 11 01:06:57 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 16 05:21:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	168308	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	315087	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	319627	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	129960	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	135834	51.111	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 102.220%			
35) Dibromofluoromethane	8.165	113	102771	53.022	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 106.040%			
50) Toluene-d8	10.565	98	398207	53.586	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 107.180%			
62) 4-Bromofluorobenzene	12.847	95	132363	49.904	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 99.800%			
Target Compounds						
16) Acetone	4.436	43	96511	88.526	ug/l	91
20) Methylene Chloride	5.289	84	8389	3.814	ug/l #	68
25) 2-Butanone	7.488	43	25277	15.292	ug/l #	85
43) Isopropyl Acetate	8.688	43	229334	35.576	ug/l #	75
51) 4-Methyl-2-Pentanone	10.447	43	22128m	6.042	ug/l	

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

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