

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091824\
 Data File : VN083998.D
 Acq On : 18 Sep 2024 21:27
 Operator : JC\MD
 Sample : P4030-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-S-08-202409

Manual Integrations
 APPROVED

Reviewed By : John Carlone 09/19/2024
 Supervised By : Mahesh Dadoda 09/19/2024

Quant Time: Sep 19 01:42:43 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	161991	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	283709	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	242209	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	106141	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	125150	51.236	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 102.480%			
35) Dibromofluoromethane	8.165	113	88950	48.974	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 97.940%			
50) Toluene-d8	10.565	98	340890	54.291	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 108.580%			
62) 4-Bromofluorobenzene	12.847	95	129976	50.173	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 100.340%			
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	21060	21.508	ug/l #	85
20) Methylene Chloride	5.271	84	6472	3.230	ug/l	96
25) 2-Butanone	7.488	43	10294	7.999	ug/l	90
39) Methylcyclohexane	9.600	83	6087	1.956	ug/l #	86
43) Isopropyl Acetate	8.688	43	272483m	60.596	ug/l	

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

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