

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111924\
 Data File : VN084952.D
 Acq On : 19 Nov 2024 17:16
 Operator : JC\MD
 Sample : P4843-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SW-WTS-01

Quant Time: Nov 20 00:32:11 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 31 18:45:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	186637	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	348139	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	291141	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	135919	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	137860	51.141	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 102.280%			
35) Dibromofluoromethane	8.165	113	111647	47.379	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 94.760%			
50) Toluene-d8	10.565	98	396481	45.675	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 91.340%			
62) 4-Bromofluorobenzene	12.847	95	155133	47.819	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 95.640%			
Target Compounds						
16) Acetone	4.465	43	40801	50.117	ug/l	97
25) 2-Butanone	7.500	43	5578	4.435	ug/l	95
30) Chloroform	7.965	83	28093	6.712	ug/l	99
40) Benzene	8.600	78	239883	22.856	ug/l	97
52) Toluene	10.629	92	8992	1.465	ug/l	98
67) Ethyl Benzene	11.965	91	312990	28.788	ug/l	99
68) m/p-Xylenes	12.070	106	21970	5.338	ug/l	99
69) o-Xylene	12.400	106	21290	5.529	ug/l	96
73) Isopropylbenzene	12.694	105	49842	5.233	ug/l	96
78) n-propylbenzene	13.035	91	28430	2.547	ug/l	99
80) 1,3,5-Trimethylbenzene	13.170	105	16882	2.138	ug/l	96
84) 1,2,4-Trimethylbenzene	13.482	105	110924	13.614	ug/l	99

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

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