

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120622\
 Data File : VN075595.D
 Acq On : 06 Dec 2022 19:25
 Operator : JC\MD
 Sample : N5876-09 50X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CHA-12-02-SIL

Quant Time: Dec 07 00:19:27 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 23 00:18:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	211811	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	341952	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	311799	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	152721	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	56652	49.827	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	99.660%		
35) Dibromofluoromethane	8.171	113	72848	54.231	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	108.460%		
50) Toluene-d8	10.571	98	160746	53.131	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	106.260%		
62) 4-Bromofluorobenzene	12.853	95	120397	53.914	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	107.820%		
Target Compounds						Qvalue
11) Tert butyl alcohol	5.547	59	16859	35.322	ug/l	100
16) Acetone	4.447	43	30079	28.363	ug/l	100
25) 2-Butanone	7.494	43	17211	11.624	ug/l	99
49) 1,4-Dioxane	9.706	88	2360	42.453	ug/l #	28
51) 4-Methyl-2-Pentanone	10.447	43	7402	2.745	ug/l #	86
52) Toluene	10.635	92	24411	4.224	ug/l	98
67) Ethyl Benzene	11.965	91	37363	3.401	ug/l	95
68) m/p-Xylenes	12.071	106	69102	15.490	ug/l	99
69) o-Xylene	12.400	106	39348	8.877	ug/l	100
80) 1,3,5-Trimethylbenzene	13.176	105	19179	2.069	ug/l	98
84) 1,2,4-Trimethylbenzene	13.488	105	100080	10.763	ug/l	99
95) Naphthalene	15.647	128	13846	1.688	ug/l	97

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

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