

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111521\
 Data File : VN069488.D
 Acq On : 15 Nov 2021 13:08
 Operator : JC/MD
 Sample : M4674-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MF-WATER-COMP

Quant Time: Nov 16 00:49:25 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 09 18:18:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1272195	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2244351	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	2204524	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	826913	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	947156	51.347	ug/l	0.00
Spiked Amount	50.000	Range 61 - 141	Recovery	= 102.700%		
35) Dibromofluoromethane	8.024	113	659504	48.995	ug/l	0.00
Spiked Amount	50.000	Range 69 - 133	Recovery	= 97.980%		
50) Toluene-d8	10.440	98	2837713	53.602	ug/l	0.00
Spiked Amount	50.000	Range 65 - 126	Recovery	= 107.200%		
62) 4-Bromofluorobenzene	12.733	95	995111	51.167	ug/l	0.00
Spiked Amount	50.000	Range 58 - 135	Recovery	= 102.340%		
Target Compounds						
						Qvalue
16) Acetone	4.296	43	64790	6.510	ug/l	89
30) Chloroform	7.815	83	124868	4.870	ug/l	100
47) Bromodichloromethane	9.759	83	23908	1.228	ug/l	94
60) Dibromochloromethane	11.240	129	7449	2.859	ug/l	97
70) Styrene	12.299	104	3560	1.590	ug/l	89
74) N-amyl acetate	12.393	43	10771	0.386	ug/l	92
87) 1,3-Dichlorobenzene	13.613	146	5862	0.225	ug/l	92
89) n-Butylbenzene	13.943	91	14059	0.326	ug/l	98
91) 1,2-Dichlorobenzene	13.994	146	6232	0.251	ug/l	97
93) 1,2,4-Trichlorobenzene	15.265	180	10764	3.889	ug/l	91
95) Naphthalene	15.509	128	42851	5.515	ug/l	100
96) 1,2,3-Trichlorobenzene	15.700	180	12200	4.088	ug/l #	87

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

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