

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120321\
 Data File : VN069714.D
 Acq On : 3 Dec 2021 13:00
 Operator : JC/MD
 Sample : M4920-03
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FDR-20211202-WC-W

Quant Time: Dec 04 05:31:26 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1374309	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2375301	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	2377128	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	864496	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	1027065	48.402	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	96.800%		
35) Dibromofluoromethane	8.024	113	704906	47.378	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	94.760%		
50) Toluene-d8	10.443	98	3019075	50.536	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	101.080%		
62) 4-Bromofluorobenzene	12.733	95	1071913	49.527	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	99.060%		
Target Compounds						Qvalue
3) Chloromethane	2.309	50	4047	0.202	ug/l	92
16) Acetone	4.293	43	276068	21.628	ug/l	96
17) Carbon Disulfide	4.527	76	7677	0.227	ug/l	97
25) 2-Butanone	7.348	43	47295	2.727	ug/l #	89
51) 4-Methyl-2-Pentanone	10.336	43	35487	1.098	ug/l	94
52) Toluene	10.505	92	14305	0.339	ug/l	98
59) 2-Hexanone	11.092	43	28271	1.157	ug/l	89
89) n-Butylbenzene	13.943	91	18519	0.393	ug/l	96
91) 1,2-Dichlorobenzene	13.991	146	8005	0.277	ug/l	90
93) 1,2,4-Trichlorobenzene	15.265	180	12908	3.736	ug/l	97
95) Naphthalene	15.509	128	61824	5.332	ug/l	96
96) 1,2,3-Trichlorobenzene	15.702	180	15560	3.506	ug/l	99

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

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