

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011022\
 Data File : VN070549.D
 Acq On : 10 Jan 2022 17:44
 Operator : JC/MD
 Sample : N1015-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 20969

Quant Time: Jan 11 06:05:21 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
 Quant Title : SW846 8260
 QLast Update : Mon Dec 27 18:47:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.083	168	1131360	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.965	114	1968145	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.741	117	1499403	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.672	152	568979	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	767688	47.194	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	94.380%		
35) Dibromofluoromethane	8.021	113	529592	44.473	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	88.940%		
50) Toluene-d8	10.438	98	2059676	42.340	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	84.680%		
62) 4-Bromofluorobenzene	12.728	95	788818	44.896	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	89.800%		
Target Compounds					Qvalue	
13) Acrolein	4.068	56	874224	397.800	ug/l #	14
16) Acetone	4.288	43	6573629	622.521	ug/l #	30
18) Methyl Acetate	4.857	43	40710	1.333	ug/l	94
25) 2-Butanone	7.340	43	2521092	189.614	ug/l	92
40) Benzene	8.459	78	20300	0.336	ug/l #	85
52) Toluene	10.502	92	25514	0.698	ug/l	95
67) Ethyl Benzene	11.841	91	14260	0.245	ug/l	96
68) m/p-Xylenes	11.945	106	22961	1.082	ug/l	98
69) o-Xylene	12.278	106	16587	0.786	ug/l	98
80) 1,3,5-Trimethylbenzene	13.058	105	44965	1.062	ug/l	98
84) 1,2,4-Trimethylbenzene	13.367	105	86504	2.122	ug/l	94
95) Naphthalene	15.504	128	55608	3.865	ug/l #	90

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

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