

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080422\
 Data File : VN073749.D
 Acq On : 04 Aug 2022 20:26
 Operator : JC\MD
 Sample : N4056-01 100X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TOTE

Quant Time: Aug 05 00:34:33 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 29 02:20:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	500917	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	775570	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	704497	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	366047	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	263837	43.220	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	86.440%
35) Dibromofluoromethane	7.951	113	253085	51.967	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.940%
50) Toluene-d8	10.380	98	841579	43.914	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	87.820%
62) 4-Bromofluorobenzene	12.674	95	323634	50.457	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.920%
Target Compounds						
					Qvalue	
16) Acetone	4.222	43	579398	176.712	ug/l	99
18) Methyl Acetate	4.787	43	8495	1.041	ug/l	96
25) 2-Butanone	7.269	43	6050	1.188	ug/l	98
37) Ethyl Acetate	7.345	43	10273	1.124	ug/l	98
40) Benzene	8.386	78	6644	0.295	ug/l #	82
52) Toluene	10.445	92	15778	1.146	ug/l	97
59) 2-Hexanone	11.033	43	6919	1.031	ug/l	85
67) Ethyl Benzene	11.786	91	10491	0.416	ug/l	95
68) m/p-Xylenes	11.898	106	12395	1.287	ug/l	97
84) 1,2,4-Trimethylbenzene	13.310	105	12857	0.593	ug/l	94
86) p-Isopropyltoluene	13.557	119	342878	16.112	ug/l	97
95) Naphthalene	15.433	128	54992	2.644	ug/l	99

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

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