

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042722\
 Data File : VN072224.D
 Acq On : 27 Apr 2022 18:29
 Operator : JC\MD
 Sample : N2575-02 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31799

Quant Time: Apr 28 06:05:53 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042622W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 26 18:50:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	271196	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.963	114	501980	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	542244	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.669	152	212855	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	208178	50.001	ug/l	0.00
Spiked Amount	50.000	Range 61 - 141	Recovery	= 100.000%		
35) Dibromofluoromethane	8.022	113	153836	47.880	ug/l	0.00
Spiked Amount	50.000	Range 69 - 133	Recovery	= 95.760%		
50) Toluene-d8	10.439	98	644711	51.066	ug/l	0.00
Spiked Amount	50.000	Range 65 - 126	Recovery	= 102.140%		
62) 4-Bromofluorobenzene	12.727	95	262889	62.735	ug/l	0.00
Spiked Amount	50.000	Range 58 - 135	Recovery	= 125.460%		
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	22675	12.632	ug/l	99
18) Methyl Acetate	4.863	43	3917	0.741	ug/l	# 67
20) Methylene Chloride	5.099	84	1985	0.577	ug/l	# 71
25) 2-Butanone	7.334	43	17740	6.206	ug/l	94
52) Toluene	10.498	92	16113	1.858	ug/l	97
59) 2-Hexanone	11.086	43	2537	0.614	ug/l	77
67) Ethyl Benzene	11.839	91	8113	0.425	ug/l	# 89
68) m/p-Xylenes	11.945	106	11276	1.589	ug/l	92
69) o-Xylene	12.275	106	5085	0.739	ug/l	100
84) 1,2,4-Trimethylbenzene	13.363	105	8414	0.555	ug/l	95
86) p-Isopropyltoluene	13.610	119	7118	0.509	ug/l	95

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

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