

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN083122\
 Data File : VN074226.D
 Acq On : 31 Aug 2022 20:06
 Operator : JC\MD
 Sample : N4444-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WH-1

Quant Time: Sep 01 00:30:13 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N083022W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 31 07:02:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	187783	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.904	114	346811	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	331237	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	162239	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	156914	53.867	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.740%	
35) Dibromofluoromethane	7.951	113	126508	56.057	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 112.120%	
50) Toluene-d8	10.381	98	407536	45.855	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 91.720%	
62) 4-Bromofluorobenzene	12.669	95	160637	59.435	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 118.880%	
Target Compounds						
11) Tert butyl alcohol	5.287	59	206207	317.112	ug/l	Qvalue 98
16) Acetone	4.222	43	1996071	1025.475	ug/l	93
25) 2-Butanone	7.263	43	194957	62.610	ug/l	# 89
51) 4-Methyl-2-Pentanone	10.269	43	13901	2.327	ug/l	# 83
52) Toluene	10.445	92	98496	14.534	ug/l	93
59) 2-Hexanone	11.028	43	23378	5.156	ug/l	81
67) Ethyl Benzene	11.786	91	5648	0.422	ug/l	# 60
68) m/p-Xylenes	11.892	106	6096	1.227	ug/l	88
69) o-Xylene	12.222	106	5003	1.010	ug/l	98
84) 1,2,4-Trimethylbenzene	13.310	105	16942	1.554	ug/l	84

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

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