

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011022\
 Data File : VN070543.D
 Acq On : 10 Jan 2022 15:13
 Operator : JC/MD
 Sample : N1012-12DL2 20X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31244DL2

Quant Time: Jan 11 05:34:41 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	8.086	168	883855	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.965	114	1493422	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1331342	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	446807	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	465261	36.611	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	73.220%
35) Dibromofluoromethane	8.021	113	325589	36.033	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	72.060%
50) Toluene-d8	10.440	98	1318558	35.721	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	71.440%
62) 4-Bromofluorobenzene	12.731	95	483764	36.286	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	72.580%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	5.372	59	57972	27.397	ug/l #	86
16) Acetone	4.285	43	2037511	246.984	ug/l	97
25) 2-Butanone	7.337	43	181576	17.481	ug/l #	87
51) 4-Methyl-2-Pentanone	10.331	43	56664	2.952	ug/l	96
52) Toluene	10.502	92	308668	11.128	ug/l	97
67) Ethyl Benzene	11.843	91	55956	1.081	ug/l	99
68) m/p-Xylenes	11.950	106	90547	4.806	ug/l	92
69) o-Xylene	12.278	106	43658	2.330	ug/l	92

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

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