

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011022\
 Data File : VN070556.D
 Acq On : 10 Jan 2022 20:41
 Operator : JC/MD
 Sample : N1023-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 HH-DRUM-COMP

Quant Time: Jan 11 06:08:01 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	1015232	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.965	114	1739928	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1572012	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.672	152	646334	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	715415	49.011	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	98.020%
35) Dibromofluoromethane	8.021	113	512195	48.653	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	97.300%
50) Toluene-d8	10.440	98	2130442	49.539	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	99.080%
62) 4-Bromofluorobenzene	12.728	95	846893	54.523	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	109.040%
Target Compounds						
					Qvalue	
16) Acetone	4.288	43	236559	24.965	ug/l	99
20) Methylene Chloride	5.098	84	88703	6.766	ug/l	91
25) 2-Butanone	7.337	43	247131	20.713	ug/l #	89
32) 1,1,1-Trichloroethane	8.016	97	30486	1.583	ug/l #	51
43) Isopropyl Acetate	8.560	43	128105	3.599	ug/l #	90
51) 4-Methyl-2-Pentanone	10.328	43	194029	8.676	ug/l	93
52) Toluene	10.502	92	68401	2.117	ug/l	97
68) m/p-Xylenes	11.948	106	40750	1.832	ug/l	99
69) o-Xylene	12.277	106	28167	1.273	ug/l	92
84) 1,2,4-Trimethylbenzene	13.366	105	50332	1.087	ug/l	88
95) Naphthalene	15.501	128	129535	6.037	ug/l	98

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

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