

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
 Data File : VN072764.D
 Acq On : 10 Jun 2022 14:57
 Operator : JC\MD
 Sample : N3160-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-2

Quant Time: Jun 11 01:54:17 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 09 05:17:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.080	168	168432	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.963	114	336344	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	282587	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	105073	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	259179	57.333	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 114.660%	
35) Dibromofluoromethane	8.022	113	186479	56.032	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 112.060%	
50) Toluene-d8	10.439	98	583264	43.926	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 87.860%	
62) 4-Bromofluorobenzene	12.727	95	219418	48.156	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 96.320%	
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	43056	19.877	ug/l	93
17) Carbon Disulfide	4.528	76	19698	3.774	ug/l	100
25) 2-Butanone	7.328	43	18895	5.423	ug/l	# 79
30) Chloroform	7.816	83	6391	1.038	ug/l	93
73) Isopropylbenzene	12.580	105	3005	0.267	ug/l	92
78) n-propylbenzene	12.921	91	4418	0.359	ug/l	90
85) sec-Butylbenzene	13.498	105	19734	1.865	ug/l	# 52
89) n-Butylbenzene	13.939	91	9413	1.342	ug/l	# 73

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

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