

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090420\
 Data File : VN063254.D
 Acq On : 4 Sep 2020 17:48
 Operator : JC/MD
 Sample : L3866-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TOTE-31230

Quant Time: Sep 07 01:51:52 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N083120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 01 04:21:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	270494	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	539127	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	544415	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	226023	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	218751	49.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.94%	
35) Dibromofluoromethane	7.55	113	166275	48.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.14%	
50) Toluene-d8	10.06	98	684430	51.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.82%	
62) 4-Bromofluorobenzene	12.38	95	243537	52.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.60%	
Target Compounds						
					Qvalue	
16) Acetone	3.78	43	58410	46.468	ug/l	98
17) Carbon Disulfide	4.02	76	7379	0.835	ug/l	98
20) Methylene Chloride	4.50	84	16210	4.135	ug/l	86
25) 2-Butanone	6.79	43	45889	23.577	ug/l	92
43) Isopropyl Acetate	8.13	43	51318	6.071	ug/l #	90
51) 4-Methyl-2-Pentanone	9.95	43	16776	3.652	ug/l	90
52) Toluene	10.13	92	70231	7.402	ug/l	99
59) 2-Hexanone	10.73	43	50695	15.705	ug/l	85

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

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