

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN070120\
 Data File : VN062333.D
 Acq On : 2 Jul 2020 9:34
 Operator : JC/MD
 Sample : L3171-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RECLAIM-SYSTEM

Quant Time: Jul 02 11:17:38 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 24 15:33:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	195797	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	408676	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	405236	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	152850	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	7.99	65	165797	60.34	ug/l	0.00
Spiked Amount			Recovery	=	120.68%	
35) Dibromofluoromethane	7.55	113	131896	50.31	ug/l	0.00
Spiked Amount			Recovery	=	100.62%	
50) Toluene-d8	10.06	98	534064	50.96	ug/l	0.00
Spiked Amount			Recovery	=	101.92%	
62) 4-Bromofluorobenzene	12.38	95	173859	49.10	ug/l	0.00
Spiked Amount			Recovery	=	98.20%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	3.78	43	157473	181.475	ug/l	96
17) Carbon Disulfide	4.01	76	10457	1.673	ug/l	97
20) Methylene Chloride	4.50	84	33124	12.972	ug/l	99
25) 2-Butanone	6.80	43	13031	9.703	ug/l	98
30) Chloroform	7.33	83	109048	23.882	ug/l	99
43) Isopropyl Acetate	8.13	43	11098	2.001	ug/l #	92
51) 4-Methyl-2-Pentanone	9.95	43	19589	6.047	ug/l	99
86) p-Isopropyltoluene	13.26	119	36673	3.859	ug/l	91

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

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