

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN063020\
 Data File : VN062267.D
 Acq On : 1 Jul 2020 3:30
 Operator : JC/MD
 Sample : L3153-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP1-A

Quant Time: Jul 01 06:05:34 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	226420	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	452184	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	443435	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	172078	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	175322	55.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.36%	
35) Dibromofluoromethane	7.55	113	137751	47.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.96%	
50) Toluene-d8	10.06	98	595551	51.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.72%	
62) 4-Bromofluorobenzene	12.38	95	208349	53.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.36%	
Target Compounds						
16) Acetone	3.78	43	71457	68.709	ug/l	98
18) Methyl Acetate	4.29	43	4539	1.797	ug/l #	74
20) Methylene Chloride	4.51	84	24231	8.126	ug/l	95
43) Isopropyl Acetate	8.13	43	38659	6.301	ug/l	98
52) Toluene	10.12	92	12001	1.420	ug/l	99
79) 2-Chlorotoluene	12.65	91	89656	9.733	ug/l	99
82) 4-Chlorotoluene	12.75	91	98533	10.344	ug/l	99
95) Naphthalene	15.11	128	3336	4.369	ug/l #	95

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

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