

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080320\
 Data File : VN062735.D
 Acq On : 3 Aug 2020 15:01
 Operator : JC/MD
 Sample : L3539-02 10X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 I-LDC-5

Quant Time: Aug 04 07:51:41 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 21 18:48:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	126472	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	221869	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	220114	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	87436	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	107688	54.81	ug/l	0.00
Spiked Amount			Recovery	=	109.62%	
35) Dibromofluoromethane	7.55	113	73031	50.11	ug/l	0.00
Spiked Amount			Recovery	=	100.22%	
50) Toluene-d8	10.06	98	291200	50.08	ug/l	0.00
Spiked Amount			Recovery	=	100.16%	
62) 4-Bromofluorobenzene	12.38	95	108657	51.11	ug/l	0.00
Spiked Amount			Recovery	=	102.22%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	3.78	43	41474	56.097	ug/l	98
17) Carbon Disulfide	4.01	76	449654	101.059	ug/l	99
25) 2-Butanone	6.79	43	6132	5.446	ug/l	91
30) Chloroform	7.33	83	2567	0.861	ug/l	93
40) Benzene	8.00	78	40788	5.697	ug/l	99
52) Toluene	10.13	92	7865	1.892	ug/l	95
67) Ethyl Benzene	11.49	91	3282	0.388	ug/l	93
68) m/p-Xylenes	11.59	106	6347	2.087	ug/l	93
69) o-Xylene	11.92	106	2356	0.816	ug/l	85
84) 1,2,4-Trimethylbenzene	13.02	105	4189	0.745	ug/l	97
95) Naphthalene	15.10	128	575827	140.710	ug/l	99

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

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