

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072820\
 Data File : VN062639.D
 Acq On : 28 Jul 2020 18:19
 Operator : JC/MD
 Sample : L3432-08
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31165-L

Quant Time: Jul 29 05:10:42 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	151826	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	266153	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	261666	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	91064	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	126493	53.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.26%	
35) Dibromofluoromethane	7.55	113	87073	49.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.62%	
50) Toluene-d8	10.06	98	347498	49.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.64%	
62) 4-Bromofluorobenzene	12.38	95	130762	51.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.54%	
Target Compounds						
16) Acetone	3.78	43	29295	33.007	ug/l	92
17) Carbon Disulfide	4.01	76	56945	10.661	ug/l	98
18) Methyl Acetate	4.29	43	2959	1.317	ug/l #	84
25) 2-Butanone	6.80	43	10407	7.699	ug/l	97
30) Chloroform	7.33	83	37926	10.597	ug/l	100
51) 4-Methyl-2-Pentanone	9.95	43	18527	6.281	ug/l	96
52) Toluene	10.12	92	5855	1.174	ug/l	83

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

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