

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092420\
 Data File : VN063660.D
 Acq On : 25 Sep 2020 5:01
 Operator : JC/MD
 Sample : L4105-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MTB-SP

Quant Time: Sep 25 06:19:04 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	193687	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	379755	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	388271	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	163951	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	175187	53.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.26%	
35) Dibromofluoromethane	7.55	113	126418	48.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.18%	
50) Toluene-d8	10.06	98	499957	50.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.16%	
62) 4-Bromofluorobenzene	12.38	95	189530	52.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.60%	
Target Compounds						
16) Acetone	3.78	43	23162	21.859	ug/l	97
18) Methyl Acetate	4.29	43	5012	1.863	ug/l #	76
20) Methylene Chloride	4.51	84	15592	5.808	ug/l	98
43) Isopropyl Acetate	8.13	43	52608	7.410	ug/l	100

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

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