

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091820\
 Data File : VN063465.D
 Acq On : 18 Sep 2020 18:13
 Operator : JC/MD
 Sample : L4059-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B33-091720-18-19

Manual Integrations
 APPROVED

MMDadoda
 9/21/2020 11:49:50 AM

Quant Time: Sep 21 02:22:55 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	240597	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	456053	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	465592	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	196028	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	198913	48.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.14%	
35) Dibromofluoromethane	7.55	113	145549	46.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.20%	
50) Toluene-d8	10.06	98	593953	50.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.06%	
62) 4-Bromofluorobenzene	12.38	95	236599	54.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.74%	
Target Compounds						
16) Acetone	3.78	43	32859	24.964	ug/l	98
18) Methyl Acetate	4.29	43	3662m	1.096	ug/l	
20) Methylene Chloride	4.51	84	14499	4.348	ug/l	96
43) Isopropyl Acetate	8.13	43	89626	10.511	ug/l	100

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

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