

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091720\
 Data File : VN063431.D
 Acq On : 17 Sep 2020 20:00
 Operator : JC/MD
 Sample : L4015-22
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 REACTOR-5W-(D-4)

Manual Integrations
 APPROVED

MMDadoda
 9/18/2020 2:45:28 PM

Quant Time: Sep 18 04:12:49 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	225511	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	423819	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	431333	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	186744	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	189803	49.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.88%	
35) Dibromofluoromethane	7.55	113	138777	47.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.60%	
50) Toluene-d8	10.06	98	557834	50.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.12%	
62) 4-Bromofluorobenzene	12.38	95	221538	54.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.56%	
Target Compounds						
16) Acetone	3.78	43	14031	11.373	ug/l	100
18) Methyl Acetate	4.30	43	5555m	1.774	ug/l	
20) Methylene Chloride	4.51	84	13906	4.449	ug/l	95
43) Isopropyl Acetate	8.13	43	81295	10.259	ug/l	99

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

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