

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091720\
 Data File : VN063430.D
 Acq On : 17 Sep 2020 19:36
 Operator : JC/MD
 Sample : L4015-21
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

MMDadoda
 9/18/2020 2:45:25 PM

Quant Time: Sep 18 04:11:09 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	233660	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	442660	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	444928	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	190829	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	198140	49.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.64%	
35) Dibromofluoromethane	7.55	113	145378	47.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.88%	
50) Toluene-d8	10.06	98	575798	49.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.94%	
62) 4-Bromofluorobenzene	12.38	95	226373	53.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.18%	
Target Compounds						
16) Acetone	3.78	43	14655	11.464	ug/l	100
18) Methyl Acetate	4.30	43	5832m	1.797	ug/l	
20) Methylene Chloride	4.51	84	13390	4.134	ug/l	88
43) Isopropyl Acetate	8.13	43	78026	9.428	ug/l	100

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

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