

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091720\
 Data File : VN063426.D
 Acq On : 17 Sep 2020 18:01
 Operator : JC/MD
 Sample : L4015-17
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 REACTOR-5W-(C-0)

Quant Time: Sep 18 04:04:54 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	237247	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	441295	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	446253	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	195991	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	192680	47.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.42%	
35) Dibromofluoromethane	7.55	113	141989	46.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.96%	
50) Toluene-d8	10.06	98	571647	49.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.52%	
62) 4-Bromofluorobenzene	12.38	95	228627	54.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.58%	
Target Compounds						
16) Acetone	3.78	43	15631	12.043	ug/l	92
18) Methyl Acetate	4.28	43	5331	1.618	ug/l #	83
20) Methylene Chloride	4.51	84	13665	4.155	ug/l	93
43) Isopropyl Acetate	8.13	43	63049	7.642	ug/l	100

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

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