

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091720\
 Data File : VN063428.D
 Acq On : 17 Sep 2020 18:49
 Operator : JC/MD
 Sample : L4015-19
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

MMDadoda
 9/18/2020 2:45:19 PM

Quant Time: Sep 18 04:08:00 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	225338	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	427034	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	426568	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	182986	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	186947	48.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.48%	
35) Dibromofluoromethane	7.55	113	139491	47.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.36%	
50) Toluene-d8	10.06	98	553847	49.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.64%	
62) 4-Bromofluorobenzene	12.38	95	223580	54.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.74%	
Target Compounds						
16) Acetone	3.78	43	17874	14.499	ug/l	99
18) Methyl Acetate	4.28	43	4715m	1.507	ug/l	
20) Methylene Chloride	4.51	84	12880	4.124	ug/l	96
43) Isopropyl Acetate	8.13	43	76529	9.585	ug/l	99

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

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