

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092319\
 Data File : VN058320.D
 Acq On : 24 Sep 2019 2:15
 Operator : JC/SP
 Sample : K4997-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-08-091919

Manual Integrations
 APPROVED

MMDadoda
 9/25/2019 12:49:47 AM

Quant Time: Sep 25 12:09:25 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091819W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 19 09:27:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	446616	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	750091	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	658813	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	263098	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	348254	55.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.98%	
35) Dibromofluoromethane	7.58	113	237065	51.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.98%	
50) Toluene-d8	10.09	98	965179	54.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.66%	
62) 4-Bromofluorobenzene	12.41	95	307519	46.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.20%	
Target Compounds						
16) Acetone	3.80	43	11114	5.264	ug/l	97
20) Methylene Chloride	4.53	84	67648	18.158	ug/l	99
43) Isopropyl Acetate	8.15	43	92373m	8.811	ug/l	

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

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