

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091119\
 Data File : VN057915.D
 Acq On : 11 Sep 2019 11:23
 Operator : JC/SP
 Sample : K4643-32
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T11-C2-(0)

Manual Integrations
 APPROVED

MMDadoda
 9/13/2019 11:46:33 PM

Quant Time: Sep 13 06:05:55 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	280229	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	482668	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	442606	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	227160	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	338034	59.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	119.12%	
35) Dibromofluoromethane	7.58	113	222418	49.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.86%	
50) Toluene-d8	10.09	98	908321	52.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.04%	
62) 4-Bromofluorobenzene	12.40	95	298920	45.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.70%	
Target Compounds						
16) Acetone	3.80	43	13381	7.192	ug/l	97
18) Methyl Acetate	4.32	43	8181m	1.649	ug/l	
43) Isopropyl Acetate	8.16	43	21584	2.219	ug/l	93

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

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