

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091418\
 Data File : VN051212.D
 Acq On : 15 Sep 2018 1:01
 Operator : MD\SY
 Sample : J4906-28
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP-PIT-4

Manual Integrations
 APPROVED

MMDadoda
 9/17/2018 1:25:38 PM

Quant Time: Sep 15 02:17:37 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091018W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 11 02:27:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	577943	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	919062	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	773593	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	244960	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	347925	54.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.20%	
35) Dibromofluoromethane	7.59	113	344203	49.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.68%	
50) Toluene-d8	10.09	98	1233932	47.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.18%	
62) 4-Bromofluorobenzene	12.40	95	318714	36.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	73.54%	
Target Compounds						
16) Acetone	3.82	43	44921	22.95	ug/l	97
18) Methyl Acetate	4.34	43	14340	2.78	ug/l	98
20) Methylene Chloride	4.53	84	39549225m	5566.76	ug/l	
28) Bromochloromethane	7.20	49	36160	6.87	ug/l	98
30) Chloroform	7.37	83	13358	1.14	ug/l	95
43) Isopropyl Acetate	8.17	43	295692	28.68	ug/l #	79
95) Naphthalene	15.13	128	13717	6.67	ug/l	98

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

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