

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081718\
 Data File : VN050753.D
 Acq On : 18 Aug 2018 5:48
 Operator : MD\SY
 Sample : J4521-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 72-11969

Manual Integrations
 APPROVED

MMDadoda
 8/20/2018 10:53:47 AM

Quant Time: Aug 18 06:35:13 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081418W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 14 08:07:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	601163	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	980298	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	853422	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	300594	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	412024	54.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.74%	
35) Dibromofluoromethane	7.59	113	382565	48.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.76%	
50) Toluene-d8	10.09	98	1347763	45.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.52%	
62) 4-Bromofluorobenzene	12.40	95	368627	37.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	75.76%	
Target Compounds						
16) Acetone	3.83	43	13334	5.63	ug/l	96
18) Methyl Acetate	4.34	43	14654	1.79	ug/l #	77
20) Methylene Chloride	4.52	84	40814436m	5813.29	ug/l	
28) Bromochloromethane	7.20	49	101590	16.58	ug/l	100
30) Chloroform	7.37	83	21281	1.56	ug/l	97
43) Isopropyl Acetate	8.17	43	231464	20.59	ug/l	98

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

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