

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080218\
 Data File : VN050293.D
 Acq On : 3 Aug 2018 4:12
 Operator : MD\SY
 Sample : J4304-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-A02

Manual Integrations
 APPROVED

MMDadoda
 8/3/2018 9:37:08 AM

Quant Time: Aug 03 05:00:50 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 26 18:10:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	684269	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1071672	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	901850	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	393772	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	458083	47.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.96%	
35) Dibromofluoromethane	7.59	113	411922	47.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.16%	
50) Toluene-d8	10.09	98	1445261	44.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.86%	
62) 4-Bromofluorobenzene	12.40	95	452660	42.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	85.48%	
Target Compounds						
					Qvalue	
16) Acetone	3.83	43	23441	6.189	ug/l	98
17) Carbon Disulfide	4.05	76	21123	0.874	ug/l #	95
18) Methyl Acetate	4.33	43	18752	0.532	ug/l	99
29) Tetrahydrofuran	7.22	42	145049	61.641	ug/l	98
39) Methylcyclohexane	9.09	83	3241	0.261	ug/l	86
73) Isopropylbenzene	12.25	105	10807	0.354	ug/l	97
78) n-propylbenzene	12.59	91	17332	0.508	ug/l	97
83) tert-Butylbenzene	12.99	119	8213	0.392	ug/l	81
84) 1,2,4-Trimethylbenzene	13.04	105	60944	2.517	ug/l	100
85) sec-Butylbenzene	13.17	105	19637	0.701	ug/l	80
86) p-Isopropyltoluene	13.29	119	8970	0.380	ug/l	86
89) n-Butylbenzene	13.62	91	8224m	0.424	ug/l	

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

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