

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN071318\
 Data File : VN049836.D
 Acq On : 14 Jul 2018 8:57
 Operator : MD\SY
 Sample : J3980-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-E-MH-S

Manual Integrations
 APPROVED

MMDadoda
 7/16/2018 11:02:51 AM

Quant Time: Jul 16 06:45:45 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N071218W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 12 15:51:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1181366	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1915228	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1628682	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	621030	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	830493	60.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	120.86%	
35) Dibromofluoromethane	7.59	113	727972	50.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.86%	
50) Toluene-d8	10.09	98	2567248	51.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.90%	
62) 4-Bromofluorobenzene	12.40	95	717294	40.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	81.28%	
Target Compounds						
8) Diethyl Ether	3.41	74	16125	2.01	ug/l	95
16) Acetone	3.82	43	30865	7.11	ug/l	95
18) Methyl Acetate	4.34	43	47907	3.49	ug/l	99
20) Methylene Chloride	4.55	84	978624	69.78	ug/l	98
30) Chloroform	7.37	83	29532	1.07	ug/l	92
43) Isopropyl Acetate	8.17	43	414437	17.65	ug/l	97
59) 2-Hexanone	10.77	43	32916	3.88	ug/l #	58
95) Naphthalene	15.14	128	63247	6.14	ug/l	98

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

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