

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN092618\
 Data File : VN051536.D
 Acq On : 26 Sep 2018 14:03
 Operator : MD\SY
 Sample : J4774-06
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OK-03-092418

Manual Integrations
 APPROVED

mmdadoda
 9/30/2018 9:55:26 AM

Quant Time: Sep 27 17:25:08 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092318W.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 24 06:41:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	801066	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1267738	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1083836	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	410001	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	538262	53.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.28%	
35) Dibromofluoromethane	7.59	113	478583	47.21	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.42%	
50) Toluene-d8	10.09	98	1767798	46.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.62%	
62) 4-Bromofluorobenzene	12.40	95	508296	41.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	82.32%	
Target Compounds						
					Qvalue	
18) Methyl Acetate	4.34	43	17653	2.310	ug/l	95
20) Methylene Chloride	4.55	84	814337	82.009	ug/l	98
30) Chloroform	7.38	83	18083	1.058	ug/l	100
43) Isopropyl Acetate	8.17	43	362374	23.742	ug/l	95
47) Bromodichloromethane	9.40	83	4377	0.299	ug/l #	93
67) Ethyl Benzene	11.51	91	14327	0.306	ug/l	97
68) m/p-Xylenes	11.62	106	18895	1.042	ug/l	94
69) o-Xylene	11.95	106	7407	0.425	ug/l	90
80) 1,3,5-Trimethylbenzene	12.73	105	7708	0.273	ug/l	94
84) 1,2,4-Trimethylbenzene	13.04	105	23349	0.821	ug/l	94
95) Naphthalene	15.14	128	19526	9.119	ug/l	98

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

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