

Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
 Data File : BM010006.D
 Acq On : 12 May 2017 04:06
 Operator : SJ/MA
 Sample : I3079-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LF001MW001-170509

Quant Time: May 12 07:49:49 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 11 19:50:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	57815	20.00	ng	-0.02
21) Naphthalene-d8	10.56	136	222667	20.00	ng	-0.02
38) Acenaphthene-d10	14.42	164	136109	20.00	ng	-0.02
63) Phenanthrene-d10	17.17	188	285979	20.00	ng	-0.03
75) Chrysene-d12	21.36	240	334840	20.00	ng	-0.03
86) Perylene-d12	23.65	264	297095	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	241110	65.26	ng	-0.01
7) Phenol-d6	6.95	99	226375	38.24	ng	-0.02
23) Nitrobenzene-d5	8.93	82	562300	92.29	ng	-0.02
41) 2,4,6-Tribromophenol	15.92	330	230548	123.77	ng	-0.02
44) 2-Fluorobiphenyl	13.03	172	919929	90.39	ng	-0.02
78) Terphenyl-d14	19.79	244	1118979	70.43	ng	-0.02
Target Compounds						
10) Phenol	6.98	94	143049	22.24	ng	Qvalue 88
12) 1,3-Dichlorobenzene	7.65	146	103215	23.05	ng	# 83
13) 1,4-Dichlorobenzene	7.80	146	280862	61.22	ng	# 92
14) 1,2-Dichlorobenzene	8.12	146	2016449	453.31	ng	# 92
17) 2-Methylphenol	8.22	107	140070	33.36	ng	# 84
20) 3+4-Methylphenols	8.55	107	705545	123.42	ng	88
22) Acetophenone	8.60	105	85726	12.59	ng	# 90
27) 2,4-Dimethylphenol	9.77	122	109900	30.07	ng	# 82
30) 1,2,4-Trichlorobenzene	10.42	180	10949	2.64	ng	# 73
31) Naphthalene	10.61	128	888587	72.49	ng	100
37) 2-Methylnaphthalene	12.22	142	78706	8.97	ng	# 90
49) Dimethylphthalate	13.88	163	26924	2.24	ng	# 96
59) Diethylphthalate	15.24	149	37342	2.96	ng	98

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

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