

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
 Data File : BG022240.D
 Acq On : 8 Jun 2016 18:10
 Operator : UM/SJ
 Sample : H3402-03 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DUP-3

Quant Time: Jun 09 06:47:17 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 06 16:12:45 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	39947	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	205188	20.00	ng	0.00
38) Acenaphthene-d10	14.75	164	122187	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	239763	20.00	ng	0.00
75) Chrysene-d12	21.77	240	208398	20.00	ng	0.00
86) Perylene-d12	25.06	264	197319	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	147929	71.60	ng	0.00
7) Phenol-d6	7.27	99	232960	71.71	ng	0.00
23) Nitrobenzene-d5	9.29	82	122599	41.66	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	75807	54.91	ng	0.00
44) 2-Fluorobiphenyl	13.37	172	358972	44.77	ng	0.00
78) Terphenyl-d14	20.08	244	378635	43.13	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.19	163	23999	2.37	ng	98
51) Acenaphthene	14.81	154	16383	2.21	ng	# 93
70) Phenanthrene	17.53	178	267766	20.56	ng	98
71) Anthracene	17.62	178	67945	5.26	ng	99
74) Fluoranthene	19.54	202	430033	28.73	ng	98
77) Pyrene	19.89	202	387657	29.92	ng	99
80) Benzo(a)anthracene	21.74	228	201595	17.25	ng	98
82) Chrysene	21.81	228	176956	15.56	ng	97
85) Indeno(1,2,3-cd)pyrene	28.83	276	97326	7.63	ng	# 87
87) Benzo(b)fluoranthene	24.01	252	212014m	18.42	ng	
88) Benzo(k)fluoranthene	24.07	252	88125m	7.78	ng	
89) Benzo(a)pyrene	24.91	252	156410	14.21	ng	# 94
91) Benzo(a,h,i)perylene	30.01	276	94934	8.80	ng	# 85

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

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