

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091415\
 Data File : BF081609.D
 Acq On : 14 Sep 2015 15:47
 Operator : UM/IZ
 Sample : G3597-03 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-3(0-5)

Quant Time: Sep 15 10:39:10 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 14 13:51:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	59702	20.00	ng	0.00
21) Naphthalene-d8	11.13	136	243441	20.00	ng	0.00
38) Acenaphthene-d10	15.26	164	107401	20.00	ng	-0.01
63) Phenanthrene-d10	18.79	188	180087	20.00	ng	0.02
75) Chrysene-d12	23.42	240	123076m	20.00	ng	0.02
86) Perylene-d12	24.85	264	99973	20.00	ng	0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	44836	11.65	ng	0.00
7) Phenol-d6	7.65	99	73114	14.35	ng	-0.01
23) Nitrobenzene-d5	9.51	82	47378	9.39	ng	-0.03
41) 2,4,6-Tribromophenol	17.17	330	12482	13.05	ng	-0.01
44) 2-Fluorobiphenyl	13.76	172	92596	11.05	ng	-0.02
78) Terphenyl-d14	22.12	244	55253m	10.45	ng	0.00

Target Compounds						Qvalue
31) Naphthalene	11.18	128	524813	40.42	ng	98
37) 2-Methylnaphthalene	12.84	142	211401	25.02	ng	# 89
45) 1,1'-Biphenyl	13.98	154	59535	6.03	ng	97
48) Acenaphthylene	14.92	152	96430	7.62	ng	98
49) Dimethylphthalate	14.82	163	24183	2.90	ng	# 97
51) Acenaphthene	15.34	154	234820	32.61	ng	# 85
54) Dibenzofuran	15.77	168	322370	30.66	ng	# 88
57) Fluorene	16.58	166	390615	48.95	ng	96
70) Phenanthrene	18.89	178	3836217	383.17	ng	96
71) Anthracene	18.98	178	907293	88.21	ng	97
72) Carbazole	19.43	167	236695	24.52	ng	# 95
74) Fluoranthene	21.49	202	2894551	267.07	ng	92
77) Pyrene	21.83	202	2676037	293.53	ng	# 91
80) Benzo(a)anthracene	23.41	228	1081326	137.96	ng	94
82) Chrysene	23.45	228	1030584m	146.68	ng	
85) Indeno(1,2,3-cd)pyrene	25.92	276	301296m	58.45	ng	
87) Benzo(b)fluoranthene	24.52	252	982510m	137.66	ng	
88) Benzo(k)fluoranthene	24.53	252	318087m	53.71	ng	
89) Benzo(a)pyrene	24.81	252	742349	122.74	ng	# 84
90) Dibenzo(a,h)anthracene	25.92	278	88765m	18.99	ng	
91) Benzo(g,h,i)perylene	26.23	276	251698	56.43	ng	# 69

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

