

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\
 Data File : BF087863.D
 Acq On : 9 Jun 2016 20:22
 Operator : UM/SJ
 Sample : H3380-03 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T5-B1(0-12)C

Quant Time: Jun 10 11:08:04 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 18:51:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	118117	20.00	ng	-0.02
21) Naphthalene-d8	8.11	136	441562	20.00	ng	-0.01
38) Acenaphthene-d10	9.85	164	175427	20.00	ng	-0.03
63) Phenanthrene-d10	11.34	188	258617	20.00	ng	-0.02
75) Chrysene-d12	13.95	240	226159	20.00	ng	-0.03
86) Perylene-d12	15.36	264	197090	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	177510	25.69	ng	-0.02
7) Phenol-d6	6.44	99	226797	27.09	ng	-0.03
23) Nitrobenzene-d5	7.38	82	124732	16.50	ng	-0.03
41) 2,4,6-Tribromophenol	10.64	330	35549	19.19	ng	-0.03
44) 2-Fluorobiphenyl	9.18	172	234480	16.54	ng	-0.03
78) Terphenyl-d14	12.91	244	157304	15.83	ng	-0.03
Target Compounds						Qvalue
31) Naphthalene	8.14	128	1854160	84.85	ng	98
37) 2-Methylnaphthalene	8.82	142	728825	50.60	ng	98
45) 1,1'-Biphenyl	9.28	154	239590	13.91	ng	97
48) Acenaphthylene	9.71	152	70027	3.88	ng	# 97
51) Acenaphthene	9.90	154	862796	76.59	ng	98
54) Dibenzofuran	10.06	168	774948	49.95	ng	# 88
57) Fluorene	10.40	166	725589	72.13	ng	100
70) Phenanthrene	11.37	178	1938601	142.85	ng	98
71) Anthracene	11.42	178	573962	41.93	ng	98
72) Carbazole	11.56	167	246206	19.22	ng	97
74) Fluoranthene	12.55	202	1358975	94.89	ng	99
77) Pvrene	12.78	202	1145361	68.25	ng	100
80) Benzo(a)anthracene	13.94	228	319060	24.76	ng	98
82) Chrysene	13.99	228	227457m	18.76	ng	
85) Indeno(1,2,3-cd)pyrene	16.72	276	27779	2.47	ng	# 100
87) Benzo(b)fluoranthene	14.96	252	184626m	13.44	ng	
88) Benzo(k)fluoranthene	14.98	252	35983m	3.47	ng	
89) Benzo(a)pyrene	15.30	252	87616	7.88	ng	98
91) Benzo(a,h,i)perylene	17.13	276	25350	2.39	ng	96

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

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