

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110415\
 Data File : BF082690.D
 Acq On : 4 Nov 2015 8:13
 Operator : UM/IZ
 Sample : G4241-02
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-117(3)

Quant Time: Nov 04 13:02:56 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF103015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 03 12:27:47 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	158605	20.00	ng	0.01
21) Naphthalene-d8	8.17	136	526520	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	222668	20.00	ng	0.01
63) Phenanthrene-d10	11.44	188	338052	20.00	ng	0.02
75) Chrysene-d12	14.09	240	387086	20.00	ng	0.03
86) Perylene-d12	15.54	264	323870	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	1267535	120.37	ng	0.01
7) Phenol-d6	6.51	99	1834177	131.11	ng	0.00
23) Nitrobenzene-d5	7.45	82	1250539	96.99	ng	0.00
41) 2,4,6-Tribromophenol	10.74	330	284982	116.95	ng	0.02
44) 2-Fluorobiphenyl	9.25	172	1245800	79.44	ng	0.00
78) Terphenyl-d14	13.03	244	991086	56.03	ng	0.02
Target Compounds						Qvalue
31) Naphthalene	8.19	128	820411	28.80	ng	99
37) 2-Methylnaphthalene	8.89	142	419226	22.70	ng	100
45) 1,1'-Biphenyl	9.36	154	105243	5.60	ng	94
48) Acenaphthylene	9.80	152	372416	15.85	ng	# 86
49) Dimethylphthalate	9.65	163	318759	17.53	ng	# 92
51) Acenaphthene	9.97	154	555561	37.25	ng	96
57) Fluorene	10.50	166	670199	41.06	ng	# 90
70) Phenanthrene	11.47	178	2228511	116.88	ng	97
71) Anthracene	11.52	178	756530m	39.41	ng	
74) Fluoranthene	12.68	202	972047	49.87	ng	92
77) Pyrene	12.90	202	1842741	64.29	ng	96
80) Benzo(a)anthracene	14.08	228	608002m	24.73	ng	
82) Chrysene	14.11	228	490903m	22.03	ng	
85) Indeno(1,2,3-cd)pyrene	16.97	276	192775	10.48	ng	98
87) Benzo(b)fluoranthene	15.13	252	471658m	21.12	ng	
88) Benzo(k)fluoranthene	15.14	252	107711m	6.33	ng	
89) Benzo(a)pyrene	15.48	252	397701	22.61	ng	# 86
90) Dibenzo(a,h)anthracene	16.98	278	36655	2.23	ng	# 1
91) Benzo(a,h,i)perylene	17.40	276	211544	13.30	ng	# 88

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

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