

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040615\
 Data File : BF078280.D
 Acq On : 7 Apr 2015 3:40
 Operator : TP/IZ
 Sample : G1745-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3D

Quant Time: Apr 07 11:43:58 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 07 04:11:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	37857	20.00	ng	0.00
21) Naphthalene-d8	8.57	136	140840	20.00	ng	0.01
38) Acenaphthene-d10	10.73	164	82470	20.00	ng	0.01
63) Phenanthrene-d10	12.56	188	146769	20.00	ng	0.01
75) Chrysene-d12	15.83	240	139918	20.00	ng	0.01
86) Perylene-d12	17.47	264	182872	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	227801	99.21	ng	0.01
7) Phenol-d6	6.58	99	296636	103.09	ng	0.00
23) Nitrobenzene-d5	7.68	82	162771	70.05	ng	0.00
41) 2,4,6-Tribromophenol	11.70	330	62177	90.91	ng	0.00
44) 2-Fluorobiphenyl	9.90	172	359003	62.22	ng	0.00
78) Terphenyl-d14	14.55	244	397899	64.26	ng	0.00

Target Compounds						Qvalue
31) Naphthalene	8.62	128	8458908	1153.93	ng	# 91
37) 2-Methylnaphthalene	9.46	142	2159293	433.86	ng	# 95
45) 1,1'-Biphenyl	10.03	154	1198680	177.69	ng	97
48) Acenaphthylene	10.54	152	681264	79.49	ng	# 96
49) Dimethylphthalate	10.42	163	12508	2.02	ng	# 99
51) Acenaphthene	10.77	154	2212203	458.45	ng	98
54) Dibenzofuran	10.98	168	114154	15.49	ng	# 95
57) Fluorene	11.41	166	1542389	258.18	ng	98
70) Phenanthrene	12.61	178	5092967	599.89	ng	93
71) Anthracene	12.66	178	1469698	175.68	ng	97
72) Carbazole	12.86	167	35031	4.47	ng	# 88
74) Fluoranthene	14.08	202	3240752	361.96	ng	92
77) Pvrene	14.35	202	4511716m	513.66	ng	
80) Benzo(a)anthracene	15.83	228	1517943	182.34	ng	# 92
82) Chrysene	15.86	228	1132618	144.81	ng	95
85) Indeno(1,2,3-cd)pyrene	18.75	276	631642m	71.17	ng	
87) Benzo(b)fluoranthene	17.07	252	1181175m	104.51	ng	
88) Benzo(k)fluoranthene	17.09	252	355856m	35.43	ng	
89) Benzo(a)pvrene	17.43	252	1328989	134.74	ng	# 94
90) Dibenzo(a,h)anthracene	18.76	278	114408	11.59	ng	97
91) Benzo(q,h,i)perylene	19.14	276	823938	83.22	ng	96

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

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