

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
 Data File : BF078343.D
 Acq On : 8 Apr 2015 19:36
 Operator : TP/IZ
 Sample : G1793-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-6DA

Quant Time: Apr 09 01:32:45 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 09 00:49:15 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.91	152	39548	20.00	ng	0.00
21) Naphthalene-d8	8.50	136	155711	20.00	ng	0.01
38) Acenaphthene-d10	10.65	164	80178	20.00	ng	0.00
63) Phenanthrene-d10	12.49	188	142016	20.00	ng	0.01
75) Chrysene-d12	15.78	240	129768	20.00	ng	0.01
86) Perylene-d12	17.57	264	149073	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	99949	41.67	ng	0.00
7) Phenol-d6	6.51	99	156121	51.94	ng	0.00
23) Nitrobenzene-d5	7.61	82	102664	39.96	ng	0.00
41) 2,4,6-Tribromophenol	11.63	330	50157	75.43	ng	0.00
44) 2-Fluorobiphenyl	9.84	172	237461	42.33	ng	0.00
78) Terphenyl-d14	14.48	244	301602	52.52	ng	0.00

Target Compounds						Qvalue
31) Naphthalene	8.53	128	4433738	547.07	ng	# 94
37) 2-Methylnaphthalene	9.38	142	1523682	276.91	ng	96
45) 1,1'-Biphenyl	9.96	154	648743	98.92	ng	99
48) Acenaphthylene	10.48	152	391398	46.97	ng	# 94
49) Dimethylphthalate	10.35	163	21478	3.57	ng	# 98
51) Acenaphthene	10.70	154	1278767	272.58	ng	99
54) Dibenzofuran	10.91	168	856043	119.47	ng	98
57) Fluorene	11.34	166	1478020	254.47	ng	98
70) Phenanthrene	12.53	178	4406142	536.36	ng	94
71) Anthracene	12.59	178	1770668	218.74	ng	96
72) Carbazole	12.78	167	259775	34.24	ng	98
74) Fluoranthene	14.00	202	2687854	310.25	ng	97
77) Pyrene	14.27	202	2713489	333.10	ng	# 92
80) Benzo(a)anthracene	15.77	228	1488579	192.80	ng	94
82) Chrysene	15.81	228	1222894m	168.58	ng	
85) Indeno(1,2,3-cd)pyrene	18.90	276	502957	61.10	ng	# 87
87) Benzo(b)fluoranthene	17.12	252	1376276m	149.38	ng	
88) Benzo(k)fluoranthene	17.14	252	274169m	33.49	ng	
89) Benzo(a)pyrene	17.52	252	1096460	136.36	ng	# 94
90) Dibenzo(a,h)anthracene	18.91	278	132243	16.43	ng	# 92
91) Benzo(g,h,i)perylene	19.27	276	464627	57.57	ng	98

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

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