

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078516.D
 Acq On : 14 Apr 2015 23:20
 Operator : TP/IZ
 Sample : G1828-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB01B

Quant Time: Apr 15 02:25:42 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 01:10:59 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	18583	20.00	ng	0.00
21) Naphthalene-d8	8.39	136	81650	20.00	ng	0.00
38) Acenaphthene-d10	10.54	164	42797	20.00	ng	0.00
63) Phenanthrene-d10	12.36	188	88557	20.00	ng	0.00
75) Chrysene-d12	15.62	240	101763	20.00	ng	-0.03
86) Perylene-d12	17.29	264	106224	20.00	ng	-0.19

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.06	112	156851	139.17	ng	0.01
7) Phenol-d6	6.41	99	208447	147.57	ng	0.00
23) Nitrobenzene-d5	7.51	82	113704	84.41	ng	0.00
41) 2,4,6-Tribromophenol	11.52	330	50596	142.55	ng	0.00
44) 2-Fluorobiphenyl	9.74	172	167174	55.84	ng	0.00
78) Terphenyl-d14	14.35	244	228622	50.77	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.43	94	6576	3.89	ng	# 70
20) 3+4-Methylphenols	7.38	107	9860	7.14	ng	97
31) Naphthalene	8.41	128	13924	3.28	ng	99
48) Acenaphthylene	10.36	152	29255	6.58	ng	98
49) Dimethylphthalate	10.25	163	17795	5.53	ng	# 97
51) Acenaphthene	10.58	154	22480	8.98	ng	97
54) Dibenzofuran	10.80	168	17424	4.56	ng	96
57) Fluorene	11.22	166	40368	13.02	ng	99
70) Phenanthrene	12.40	178	452853	88.40	ng	100
71) Anthracene	12.46	178	157100	31.12	ng	98
72) Carbazole	12.67	167	25102	5.31	ng	95
74) Fluoranthene	13.86	202	607747	112.50	ng	99
77) Pyrene	14.14	202	557757m	87.31	ng	
80) Benzo(a)anthracene	15.61	228	303699	50.16	ng	96
82) Chrysene	15.66	228	239428	42.09	ng	97
85) Indeno(1,2,3-cd)pyrene	18.85	276	107110	16.59	ng	# 90
87) Benzo(b)fluoranthene	16.87	252	316629m	48.23	ng	
88) Benzo(k)fluoranthene	16.90	252	71812m	12.31	ng	
89) Benzo(a)pyrene	17.23	252	233485	40.75	ng	97
90) Dibenzo(a,h)anthracene	18.51	278	31926m	5.57	ng	
91) Benzo(g,h,i)perylene	18.85	276	112451m	19.55	ng	

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

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