

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078505.D
 Acq On : 14 Apr 2015 18:05
 Operator : TP/IZ
 Sample : G1828-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB05A

Quant Time: Apr 15 10:44:56 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	20523	20.00	ng	0.00
21) Naphthalene-d8	8.38	136	83844	20.00	ng	-0.01
38) Acenaphthene-d10	10.54	164	41984	20.00	ng	0.00
63) Phenanthrene-d10	12.37	188	82175	20.00	ng	0.00
75) Chrysene-d12	15.66	240	93841	20.00	ng	0.00
86) Perylene-d12	17.45	264	93094	20.00	ng	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.05	112	125205	100.59	ng	0.00
7) Phenol-d6	6.41	99	172027	110.28	ng	0.00
23) Nitrobenzene-d5	7.51	82	98024	70.86	ng	0.00
41) 2,4,6-Tribromophenol	11.52	330	41597	119.46	ng	0.00
44) 2-Fluorobiphenyl	9.74	172	195675	66.62	ng	0.00
78) Terphenyl-d14	14.37	244	237470	57.18	ng	0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.42	94	5685	3.04	ng	# 54
31) Naphthalene	8.41	128	26033	5.97	ng	98
37) 2-Methylnaphthalene	9.27	142	35164	11.87	ng	99
45) 1,1'-Biphenyl	9.85	154	11759	3.42	ng	97
51) Acenaphthene	10.58	154	14920	6.07	ng	96
54) Dibenzofuran	10.80	168	19718	5.26	ng	95
57) Fluorene	11.22	166	16934	5.57	ng	# 87
70) Phenanthrene	12.40	178	225595	47.46	ng	98
71) Anthracene	12.46	178	77143m	16.47	ng	
72) Carbazole	12.67	167	33159	7.55	ng	# 85
74) Fluoranthene	13.86	202	353329	70.48	ng	97
77) Pyrene	14.14	202	353084	59.94	ng	# 93
80) Benzo(a)anthracene	15.65	228	213800	38.29	ng	95
82) Chrysene	15.68	228	194547m	37.09	ng	
85) Indeno(1,2,3-cd)pyrene	18.74	276	197378	33.16	ng	# 86
87) Benzo(b)fluoranthene	16.99	252	351658m	61.12	ng	
88) Benzo(k)fluoranthene	17.03	252	85285m	16.68	ng	
89) Benzo(a)pyrene	17.39	252	235641	46.93	ng	# 92
90) Dibenzo(a,h)anthracene	18.75	278	44964m	8.94	ng	
91) Benzo(g,h,i)perylene	19.10	276	214686	42.59	ng	96

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

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