

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090716\
 Data File : BF089999.D
 Acq On : 7 Sep 2016 12:31
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Sep 08 10:23:12 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 02 13:29:08 2016
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	119	-0.02
2	1,4-Dioxane	0.592	0.553	6.6	118	-0.05
3	Pyridine	1.579	1.469	7.0	106	-0.05
4	n-Nitrosodimethylamine	0.779	0.754	3.2	111	-0.05
5 S	2-Fluorophenol	1.250	1.233	1.4	112	-0.02
6	Aniline	2.072	1.926	7.0	116	-0.01
7 S	Phenol-d6	1.518	1.459	3.9	113	-0.01
8	2-Chlorophenol	1.356	1.315	3.0	114	-0.02
9	Benzaldehyde	0.980	0.932	4.9	112	-0.02
10 C	Phenol	1.766	1.743	1.3	121	-0.01
11	bis(2-Chloroethyl)ether	1.336	1.338	-0.1	126	-0.02
12	1,3-Dichlorobenzene	1.514	1.540	-1.7	124	-0.02
13 C	1,4-Dichlorobenzene	1.549	1.595	-3.0	127	-0.02
14	1,2-Dichlorobenzene	1.389	1.362	1.9	113	-0.02
15	Benzyl Alcohol	0.951	0.960	-0.9	119	-0.01
16	2,2'-oxybis(1-Chloropropane	2.547	2.425	4.8	110	-0.02
17	2-Methylphenol	1.083	1.058	2.3	116	-0.01
18	Hexachloroethane	0.531	0.523	1.5	117	-0.02
19 P	n-Nitroso-di-n-propylamine	0.989	0.882	10.8	101	-0.02
20	3+4-Methylphenols	1.312	1.347	-2.7	138	-0.02
21 I	Naphthalene-d8	1.000	1.000	0.0	125	-0.02
22	Acetophenone	0.449	0.414	7.8	118	-0.02
23 S	Nitrobenzene-d5	0.345	0.317	8.1	115	-0.02
24	Nitrobenzene	0.367	0.343	6.5	125	-0.02
25	Isophorone	0.701	0.635	9.4	116	-0.02
26 C	2-Nitrophenol	0.175	0.189	-8.0	135	-0.01
27	2,4-Dimethylphenol	0.324	0.292	9.9	124	-0.02
28	bis(2-Chloroethoxy)methane	0.423	0.409	3.3	117	-0.01
29 C	2,4-Dichlorophenol	0.291	0.282	3.1	128	-0.02
30	1,2,4-Trichlorobenzene	0.318	0.305	4.1	117	-0.02
31	Naphthalene	0.996	0.931	6.5	127	-0.02
32	Benzoic acid	0.218	0.244	-11.9	141	0.00
33	4-Chloroaniline	0.402	0.408	-1.5	125	-0.01
34 C	Hexachlorobutadiene	0.186	0.188	-1.1	130	-0.02
35	Caprolactam	0.092	0.091	1.1	118	-0.01
36 C	4-Chloro-3-methylphenol	0.308	0.310	-0.6	126	-0.01
37	2-Methylnaphthalene	0.658	0.588	10.6	110	-0.02
38 I	Acenaphthene-d10	1.000	1.000	0.0	120	-0.01
39	1,2,4,5-Tetrachlorobenzene	0.586	0.607	-3.6	141	-0.01
40 P	Hexachlorocyclopentadiene	0.300	0.335	-11.7	131	-0.02
41 S	2,4,6-Tribromophenol	0.197	0.188	4.6	114	-0.02
42 C	2,4,6-Trichlorophenol	0.406	0.426	-4.9	119	-0.01
43	2,4,5-Trichlorophenol	0.390	0.408	-4.6	136	-0.02
44 S	2-Fluorobiphenyl	1.218	1.145	6.0	117	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	1,1'-Biphenyl	1.610	1.497	7.0	124	-0.02
46	2-Chloronaphthalene	1.139	1.171	-2.8	129	-0.01
47	2-Nitroaniline	0.372	0.369	0.8	123	-0.01
48	Acenaphthylene	1.884	1.815	3.7	119	-0.01
49	Dimethylphthalate	1.446	1.284	11.2	121	-0.02
50	2,6-Dinitrotoluene	0.322	0.312	3.1	120	-0.01
51 C	Acenaphthene	1.193	1.220	-2.3	129	-0.01
52	3-Nitroaniline	0.367	0.387	-5.4	141	-0.01
53 P	2,4-Dinitrophenol	0.130	0.172	-32.3#	144	-0.01
54	Dibenzofuran	1.599	1.514	5.3	117	-0.01
55 P	4-Nitrophenol	0.277	0.301	-8.7	121	0.00
56	2,4-Dinitrotoluene	0.408	0.417	-2.2	110	-0.01
57	Fluorene	1.169	1.031	11.8	102	-0.02
58	2,3,4,6-Tetrachlorophenol	0.334	0.340	-1.8	127	-0.02
59	Diethylphthalate	1.360	1.307	3.9	123	-0.01
60	4-Chlorophenyl-phenylether	0.550	0.488	11.3	106	-0.02
61	4-Nitroaniline	0.360	0.346	3.9	106	-0.01
62	Azobenzene	1.358	1.291	4.9	123	-0.02
63 I	Phenanthrene-d10	1.000	1.000	0.0	115	-0.02
64	4,6-Dinitro-2-methylphenol	0.128	0.128	0.0	126	-0.02
65 c	n-Nitrosodiphenylamine	0.601	0.622	-3.5	127	-0.01
66	4-Bromophenyl-phenylether	0.201	0.207	-3.0	123	-0.01
67	Hexachlorobenzene	0.230	0.234	-1.7	132	-0.02
68	Atrazine	0.201	0.191	5.0	119	-0.02
69 C	Pentachlorophenol	0.124	0.148	-19.4	124	-0.01
70	Phenanthrene	0.996	0.940	5.6	111	-0.02
71	Anthracene	1.010	0.989	2.1	128	-0.02
72	Carbazole	0.947	0.972	-2.6	124	-0.01
73	Di-n-butylphthalate	1.104	1.093	1.0	122	-0.02
74 C	Fluoranthene	1.044	0.915	12.4	100	-0.02
75 I	Chrysene-d12	1.000	1.000	0.0	111	-0.02
76	Benzidine	0.758	0.803	-5.9	117	-0.01
77	Pyrene	1.410	1.365	3.2	118	-0.02
78 S	Terphenyl-d14	0.810	0.800	1.2	130	-0.02
79	Butylbenzylphthalate	0.569	0.601	-5.6	120	-0.01
80	Benzo(a)anthracene	1.225	1.189	2.9	110	-0.02
81	3,3'-Dichlorobenzidine	0.442	0.485	-9.7	125	-0.02
82	Chrysene	1.144	1.078	5.8	110	-0.02
83	Bis(2-ethylhexyl)phthalate	0.787	0.793	-0.8	112	-0.02
84 c	Di-n-octyl phthalate	1.270	1.368	-7.7	111	-0.02
85	Indeno(1,2,3-cd)pyrene	0.846	0.865	-2.2	113	-0.03
86 I	Perylene-d12	1.000	1.000	0.0	112	-0.02
87	Benzo(b)fluoranthene	1.253	1.300	-3.8	102	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.063	0.949	10.7	118	-0.02
89 C	Benzo(a)pyrene	1.073	1.095	-2.1	112	-0.02
90	Dibenzo(a,h)anthracene	0.787	0.797	-1.3	113	-0.05
91	Benzo(g,h,i)perylene	0.792	0.803	-1.4	115	-0.05

(#) = Out of Range

SPCC's out = 0 CCC's out = 0