

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
 Data File : BF087477.D
 Acq On : 28 May 2016 13:38
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: May 30 03:18:15 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 25 16:26:52 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	98	-0.05
2	1,4-Dioxane	0.601	0.552	8.2	92	-0.06
3	Pyridine	1.313	1.221	7.0	85	-0.07
4	n-Nitrosodimethylamine	1.012	1.132	-11.9	106	-0.07
5 S	2-Fluorophenol	1.138	1.258	-10.5	102	-0.06
6	Aniline	1.990	2.140	-7.5	101	-0.06
7 S	Phenol-d6	1.538	1.620	-5.3	103	-0.03
8	2-Chlorophenol	1.419	1.456	-2.6	100	-0.05
9	Benzaldehyde	0.849	0.876	-3.2	110	-0.06
10 C	Phenol	1.679	1.728	-2.9	109	-0.03
11	bis(2-Chloroethyl)ether	1.359	1.363	-0.3	99	-0.05
12	1,3-Dichlorobenzene	1.548	1.543	0.3	99	-0.05
13 C	1,4-Dichlorobenzene	1.589	1.585	0.3	99	-0.06
14	1,2-Dichlorobenzene	1.470	1.465	0.3	100	-0.06
15	Benzyl Alcohol	1.121	1.276	-13.8	104	-0.05
16	2,2'-oxybis(1-Chloropropane	3.058	3.022	1.2	99	-0.05
17	2-Methylphenol	1.138	1.180	-3.7	102	-0.05
18	Hexachloroethane	0.593	0.603	-1.7	100	-0.06
19 P	n-Nitroso-di-n-propylamine	1.147	1.185	-3.3	103	-0.05
20	3+4-Methylphenols	1.375	1.535	-11.6	104	-0.05
21 I	Naphthalene-d8	1.000	1.000	0.0	100	-0.05
22	Acetophenone	0.478	0.503	-5.2	104	-0.05
23 S	Nitrobenzene-d5	0.397	0.422	-6.3	103	-0.05
24	Nitrobenzene	0.419	0.442	-5.5	101	-0.05
25	Isophorone	0.712	0.737	-3.5	103	-0.06
26 C	2-Nitrophenol	0.171	0.184	-7.6	100	-0.06
27	2,4-Dimethylphenol	0.297	0.301	-1.3	99	-0.05
28	bis(2-Chloroethoxy)methane	0.401	0.394	1.7	99	-0.05
29 C	2,4-Dichlorophenol	0.268	0.281	-4.9	100	-0.05
30	1,2,4-Trichlorobenzene	0.307	0.302	1.6	98	-0.06
31	Naphthalene	1.002	1.013	-1.1	103	-0.05
32	Benzoic acid	0.141	0.132	6.4	86	-0.02
33	4-Chloroaniline	0.369	0.385	-4.3	101	-0.06
34 C	Hexachlorobutadiene	0.186	0.186	0.0	99	-0.06
35	Caprolactam	0.078	0.084	-7.7	103	-0.02
36 C	4-Chloro-3-methylphenol	0.303	0.327	-7.9	103	-0.03
37	2-Methylnaphthalene	0.626	0.603	3.7	97	-0.06
38 I	Acenaphthene-d10	1.000	1.000	0.0	102	-0.06
39	1,2,4,5-Tetrachlorobenzene	0.640	0.615	3.9	98	-0.06
40 P	Hexachlorocyclopentadiene	0.206	0.164	20.4	71	-0.05
41 S	2,4,6-Tribromophenol	0.192	0.200	-4.2	102	-0.05
42 C	2,4,6-Trichlorophenol	0.370	0.376	-1.6	99	-0.05
43	2,4,5-Trichlorophenol	0.376	0.381	-1.3	101	-0.05
44 S	2-Fluorobiphenyl	1.419	1.370	3.5	101	-0.05

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	1,1'-Biphenyl	1.684	1.679	0.3	105	-0.05
46	2-Chloronaphthalene	1.288	1.259	2.3	101	-0.05
47	2-Nitroaniline	0.464	0.514	-10.8	109	-0.05
48	Acenaphthylene	2.040	1.937	5.0	98	-0.06
49	Dimethylphthalate	1.602	1.581	1.3	101	-0.05
50	2,6-Dinitrotoluene	0.323	0.303	6.2	95	-0.05
51 C	Acenaphthene	1.254	1.199	4.4	102	-0.05
52	3-Nitroaniline	0.331	0.357	-7.9	108	-0.05
53 P	2,4-Dinitrophenol	0.138	0.116	15.9	73	-0.02
54	Dibenzofuran	1.697	1.654	2.5	102	-0.05
55 P	4-Nitrophenol	0.158	0.166	-5.1	105	-0.02
56	2,4-Dinitrotoluene	0.431	0.455	-5.6	103	-0.03
57	Fluorene	1.472	1.455	1.2	102	-0.06
58	2,3,4,6-Tetrachlorophenol	0.289	0.280	3.1	93	-0.05
59	Diethylphthalate	1.558	1.509	3.1	101	-0.05
60	4-Chlorophenyl-phenylether	0.718	0.699	2.6	101	-0.05
61	4-Nitroaniline	0.309	0.324	-4.9	96	-0.03
62	Azobenzene	1.616	1.681	-4.0	102	-0.06
63 I	Phenanthrene-d10	1.000	1.000	0.0	104	-0.05
64	4,6-Dinitro-2-methylphenol	0.127	0.110	13.4	81	-0.03
65 c	n-Nitrosodiphenylamine	0.647	0.617	4.6	97	-0.05
66	4-Bromophenyl-phenylether	0.219	0.214	2.3	100	-0.06
67	Hexachlorobenzene	0.229	0.228	0.4	102	-0.05
68	Atrazine	0.206	0.155	24.8	77	-0.05
69 C	Pentachlorophenol	0.061	0.072	-18.0	108	-0.05
70	Phenanthrene	1.108	1.108	0.0	105	-0.05
71	Anthracene	1.119	1.081	3.4	101	-0.05
72	Carbazole	0.955	0.972	-1.8	105	-0.03
73	Di-n-butylphthalate	1.234	1.162	5.8	93	-0.05
74 C	Fluoranthene	1.111	1.055	5.0	97	-0.05
75 I	Chrysene-d12	1.000	1.000	0.0	91	-0.05
76	Benzidine	0.515	0.708	-37.5#	113	-0.05
77	Pyrene	1.440	1.583	-9.9	99	-0.05
78 S	Terphenyl-d14	0.941	1.069	-13.6	101	-0.05
79	Butylbenzylphthalate	0.638	0.674	-5.6	89	-0.05
80	Benzo(a)anthracene	1.138	1.134	0.4	90	-0.03
81	3,3'-Dichlorobenzidine	0.628	0.613	2.4	91	-0.05
82	Chrysene	1.129	1.080	4.3	89	-0.05
83	Bis(2-ethylhexyl)phthalate	0.932	0.897	3.8	85	-0.05
84 c	Di-n-octyl phthalate	1.421	1.264	11.0	75	-0.05
85	Indeno(1,2,3-cd)pyrene	0.905	1.075	-18.8	106	-0.03
86 I	Perylene-d12	1.000	1.000	0.0	85	-0.03
87	Benzo(b)fluoranthene	1.274	1.301	-2.1	81	-0.03

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88	Benzo(k)fluoranthene	1.097	0.980	10.7	90	-0.03
89 C	Benzo(a)pyrene	1.102	1.099	0.3	87	-0.03
90	Dibenzo(a,h)anthracene	0.940	1.157	-23.1	104	-0.03
91	Benzo(g,h,i)perylene	0.927	1.181	-27.4#	107	-0.05

(#) = Out of Range

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