

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050916\
 Data File : BF086828.D
 Acq On : 9 May 2016 17:50
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: May 10 02:29:58 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 09 16:04:56 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	104	0.00
2	1,4-Dioxane	0.580	0.565	2.6	101	0.00
3	Pyridine	1.417	1.494	-5.4	109	-0.01
4	n-Nitrosodimethylamine	1.028	1.056	-2.7	101	-0.01
5 S	2-Fluorophenol	1.181	1.198	-1.4	104	0.00
6	Aniline	1.909	1.864	2.4	102	0.00
7 S	Phenol-d6	1.578	1.540	2.4	101	-0.01
8	2-Chlorophenol	1.425	1.365	4.2	99	0.00
9	Benzaldehyde	0.914	0.842	7.9	99	0.00
10 C	Phenol	1.876	1.661	11.5	89	0.00
11	bis(2-Chloroethyl)ether	1.463	1.483	-1.4	99	0.00
12	1,3-Dichlorobenzene	1.542	1.412	8.4	96	-0.01
13 C	1,4-Dichlorobenzene	1.573	1.414	10.1	99	0.00
14	1,2-Dichlorobenzene	1.371	1.375	-0.3	100	0.00
15	Benzyl Alcohol	1.090	1.065	2.3	100	0.00
16	2,2'-oxybis(1-Chloropropane	3.180	2.895	9.0	100	0.00
17	2-Methylphenol	1.134	1.129	0.4	100	0.00
18	Hexachloroethane	0.592	0.552	6.8	98	-0.01
19 P	n-Nitroso-di-n-propylamine	1.103	1.126	-2.1	98	0.00
20	3+4-Methylphenols	1.481	1.441	2.7	97	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
22	Acetophenone	0.472	0.440	6.8	100	0.00
23 S	Nitrobenzene-d5	0.398	0.375	5.8	100	0.00
24	Nitrobenzene	0.410	0.401	2.2	103	0.00
25	Isophorone	0.739	0.682	7.7	103	0.00
26 C	2-Nitrophenol	0.180	0.178	1.1	99	0.00
27	2,4-Dimethylphenol	0.307	0.265	13.7	99	0.00
28	bis(2-Chloroethoxy)methane	0.399	0.396	0.8	102	0.00
29 C	2,4-Dichlorophenol	0.270	0.250	7.4	99	-0.01
30	1,2,4-Trichlorobenzene	0.302	0.292	3.3	103	0.00
31	Naphthalene	1.002	0.918	8.4	99	0.00
32	Benzoic acid	0.180	0.189	-5.0	112	0.00
33	4-Chloroaniline	0.389	0.378	2.8	101	0.00
34 C	Hexachlorobutadiene	0.175	0.165	5.7	100	0.00
35	Caprolactam	0.092	0.083	9.8	101	0.00
36 C	4-Chloro-3-methylphenol	0.327	0.300	8.3	97	0.00
37	2-Methylnaphthalene	0.586	0.579	1.2	102	0.00
38 I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
39	1,2,4,5-Tetrachlorobenzene	0.582	0.535	8.1	102	0.00
40 P	Hexachlorocyclopentadiene	0.276	0.257	6.9	98	0.00
41 S	2,4,6-Tribromophenol	0.212	0.200	5.7	106	0.00
42 C	2,4,6-Trichlorophenol	0.389	0.349	10.3	101	0.00
43	2,4,5-Trichlorophenol	0.402	0.378	6.0	102	0.00
44 S	2-Fluorobiphenyl	1.190	1.013	14.9	103	0.00

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050916\
 Data File : BF086828.D
 Acq On : 9 May 2016 17:50
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: May 10 02:29:58 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 09 16:04:56 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
45	1,1'-Biphenyl	1.591	1.565	1.6	103	0.00
46	2-Chloronaphthalene	1.246	1.210	2.9	101	0.00
47	2-Nitroaniline	0.456	0.414	9.2	104	0.00
48	Acenaphthylene	1.833	1.833	0.0	103	0.00
49	Dimethylphthalate	1.526	1.366	10.5	104	0.00
50	2,6-Dinitrotoluene	0.321	0.289	10.0	103	0.00
51 C	Acenaphthene	1.140	1.162	-1.9	103	0.00
52	3-Nitroaniline	0.338	0.312	7.7	102	0.00
53 P	2,4-Dinitrophenol	0.134	0.135	-0.7	106	0.00
54	Dibenzofuran	1.464	1.428	2.5	100	0.00
55 P	4-Nitrophenol	0.201	0.208	-3.5	105	0.00
56	2,4-Dinitrotoluene	0.397	0.383	3.5	103	0.00
57	Fluorene	1.248	1.238	0.8	103	0.00
58	2,3,4,6-Tetrachlorophenol	0.303	0.298	1.7	101	0.00
59	Diethylphthalate	1.487	1.442	3.0	104	0.00
60	4-Chlorophenyl-phenylether	0.635	0.518	18.4	102	0.00
61	4-Nitroaniline	0.325	0.335	-3.1	104	0.00
62	Azobenzene	1.604	1.506	6.1	104	0.00
63 I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
64	4,6-Dinitro-2-methylphenol	0.124	0.118	4.8	106	0.00
65 c	n-Nitrosodiphenylamine	0.614	0.583	5.0	108	0.00
66	4-Bromophenyl-phenylether	0.203	0.205	-1.0	104	0.00
67	Hexachlorobenzene	0.237	0.210	11.4	106	0.00
68	Atrazine	0.205	0.199	2.9	97	0.00
69 C	Pentachlorophenol	0.110	0.117	-6.4	106	0.00
70	Phenanthrene	1.067	0.907	15.0	101	0.00
71	Anthracene	1.092	1.092	0.0	105	0.00
72	Carbazole	0.938	0.916	2.3	109	0.00
73	Di-n-butylphthalate	1.146	1.104	3.7	104	0.00
74 C	Fluoranthene	1.098	1.064	3.1	104	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	116	0.00
76	Benzidine	0.510	0.594	-16.5	124	0.00
77	Pyrene	1.531	1.301	15.0	103	0.00
78 S	Terphenyl-d14	0.943	0.871	7.6	108	0.00
79	Butylbenzylphthalate	0.648	0.582	10.2	108	0.00
80	Benzo(a)anthracene	1.123	1.023	8.9	110	0.00
81	3,3'-Dichlorobenzidine	0.713	0.717	-0.6	111	0.00
82	Chrysene	1.142	0.984	13.8	112	0.00
83	Bis(2-ethylhexyl)phthalate	0.779	0.700	10.1	106	0.00
84 c	Di-n-octyl phthalate	1.290	1.192	7.6	109	0.00
85	Indeno(1,2,3-cd)pyrene	0.950	0.821	13.6	100	0.00
86 I	Perylene-d12	1.000	1.000	0.0	114	0.00
87	Benzo(b)fluoranthene	1.300	1.190	8.5	114	0.00

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050916\
Data File : BF086828.D
Acq On : 9 May 2016 17:50
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: May 10 02:29:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 09 16:04:56 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)
88	Benzo(k)fluoranthene	1.064	1.031	3.1	107	0.00
89 C	Benzo(a)pyrene	1.121	1.038	7.4	108	0.00
90	Dibenzo(a,h)anthracene	0.945	0.870	7.9	101	0.00
91	Benzo(a,h,i)perylene	0.960	0.887	7.6	101	0.00

(#) = Out of Range

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