

Data Path : U:\HPCHEM1\BNA F\DATA\BF082117\
 Data File : BF097909.D
 Acq On : 21 Aug 2017 17:02
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 22 04:59:30 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 21 15:59:04 2017
 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	0.733	0.723	1.4	103	0.00
3	Pyridine	2.027	2.045	-0.9	105	0.01
4	n-Nitrosodimethylamine	0.780	0.772	1.0	99	0.00
5 S	2-Fluorophenol	1.315	1.289	2.0	102	0.00
6	Aniline	2.510	2.450	2.4	102	0.00
7 S	Phenol-d6	1.787	1.799	-0.7	105	0.00
8	2-Chlorophenol	1.387	1.405	-1.3	104	0.00
9	Benzaldehyde	1.132	1.078	4.8	97	0.00
10 C	Phenol	2.089	2.122	-1.6	102	0.00
11	bis(2-Chloroethyl)ether	1.577	1.592	-1.0	105	0.00
12	1,3-Dichlorobenzene	1.492	1.473	1.3	103	0.00
13 C	1,4-Dichlorobenzene	1.517	1.491	1.7	103	0.00
14	1,2-Dichlorobenzene	1.399	1.371	2.0	102	0.00
15	Benzyl Alcohol	1.338	1.311	2.0	100	0.00
16	2,2'-oxybis(1-Chloropropane	2.122	2.061	2.9	100	0.00
17	2-Methylphenol	1.222	1.231	-0.7	104	0.00
18	Hexachloroethane	0.595	0.598	-0.5	102	0.00
19 P	n-Nitroso-di-n-propylamine	1.223	1.171	4.3	99	0.00
20	3+4-Methylphenols	1.493	1.431	4.2	99	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
22	Acetophenone	0.528	0.484	8.3	99	0.00
23 S	Nitrobenzene-d5	0.368	0.405	-10.1	113	0.00
24	Nitrobenzene	0.410	0.448	-9.3	108	0.00
25	Isophorone	0.766	0.764	0.3	105	0.00
26 C	2-Nitrophenol	0.137	0.177	-29.2#	131	0.00
27	2,4-Dimethylphenol	0.319	0.317	0.6	106	0.00
28	bis(2-Chloroethoxy)methane	0.503	0.496	1.4	104	0.00
29 C	2,4-Dichlorophenol	0.265	0.271	-2.3	106	0.00
30	1,2,4-Trichlorobenzene	0.294	0.288	2.0	104	0.00
31	Naphthalene	1.038	1.007	3.0	104	0.00
32	Benzoic acid	0.212	0.243	-14.6	122	0.00
33	4-Chloroaniline	0.438	0.435	0.7	106	0.00
34 C	Hexachlorobutadiene	0.172	0.167	2.9	103	0.00
35	Caprolactam	0.090	0.100	-11.1	113	0.00
36 C	4-Chloro-3-methylphenol	0.341	0.347	-1.8	105	0.00
37	2-Methylnaphthalene	0.637	0.626	1.7	104	0.00
38 I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
39	1,2,4,5-Tetrachlorobenzene	0.614	0.606	1.3	107	0.00
40 P	Hexachlorocyclopentadiene	0.295	0.303	-2.7	104	0.00
41 S	2,4,6-Tribromophenol	0.174	0.185	-6.3	110	0.00
42 C	2,4,6-Trichlorophenol	0.412	0.425	-3.2	108	0.00
43	2,4,5-Trichlorophenol	0.409	0.434	-6.1	110	0.00
44 S	2-Fluorobiphenyl	1.361	1.248	8.3	102	0.00

Data Path : U:\HPCHEM1\BNA F\DATA\BF082117\
 Data File : BF097909.D
 Acq On : 21 Aug 2017 17:02
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 22 04:59:30 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 21 15:59:04 2017
 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)
45	1,1'-Biphenyl	1.824	1.761	3.5	105	0.00
46	2-Chloronaphthalene	1.348	1.292	4.2	105	0.00
47	2-Nitroaniline	0.452	0.543	-20.1	122	0.00
48	Acenaphthylene	2.116	2.065	2.4	105	0.00
49	Dimethylphthalate	1.462	1.474	-0.8	108	0.00
50	2,6-Dinitrotoluene	0.292	0.324	-11.0	115	0.00
51 C	Acenaphthene	1.335	1.254	6.1	102	0.00
52	3-Nitroaniline	0.376	0.421	-12.0	117	0.00
53 P	2,4-Dinitrophenol	0.106	0.164	-54.7#	161#	0.00
54	Dibenzofuran	1.789	1.745	2.5	106	0.00
55 P	4-Nitrophenol	0.300	0.323	-7.7	110	0.00
56	2,4-Dinitrotoluene	0.366	0.417	-13.9	115	0.00
57	Fluorene	1.383	1.329	3.9	106	0.00
58	2,3,4,6-Tetrachlorophenol	0.317	0.338	-6.6	112	0.00
59	Diethylphthalate	1.529	1.504	1.6	105	0.00
60	4-Chlorophenyl-phenylether	0.642	0.618	3.7	105	0.00
61	4-Nitroaniline	0.358	0.410	-14.5	119	0.00
62	Azobenzene	1.816	1.762	3.0	104	0.00
63 I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
64	4,6-Dinitro-2-methylphenol	0.083	0.114	-37.3#	148	0.00
65 c	n-Nitrosodiphenylamine	0.681	0.653	4.1	107	0.00
66	4-Bromophenyl-phenylether	0.208	0.203	2.4	107	0.00
67	Hexachlorobenzene	0.212	0.206	2.8	107	0.00
68	Atrazine	0.181	0.176	2.8	107	0.00
69 C	Pentachlorophenol	0.123	0.130	-5.7	109	0.00
70	Phenanthrene	1.066	1.011	5.2	106	0.00
71	Anthracene	1.081	1.039	3.9	108	0.00
72	Carbazole	1.026	1.005	2.0	110	0.00
73	Di-n-butylphthalate	1.182	1.208	-2.2	111	0.00
74 C	Fluoranthene	1.112	1.102	0.9	110	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	127	0.00
76	Benzidine	0.656	0.574	12.5	118	0.00
77	Pyrene	1.636	1.440	12.0	112	0.00
78 S	Terphenyl-d14	0.959	0.818	14.7	110	0.00
79	Butylbenzylphthalate	0.627	0.647	-3.2	127	0.00
80	Benzo(a)anthracene	1.199	1.062	11.4	113	0.00
81	3,3'-Dichlorobenzidine	0.404	0.416	-3.0	124	0.00
82	Chrysene	1.192	1.088	8.7	118	0.00
83	Bis(2-ethylhexyl)phthalate	0.695	0.742	-6.8	125	0.00
84 c	Di-n-octyl phthalate	1.209	1.422	-17.6	141	0.00
85	Indeno(1,2,3-cd)pyrene	0.877	0.723	17.6	108	0.01
86 I	Perylene-d12	1.000	1.000	0.0	124	0.00
87	Benzo(b)fluoranthene	1.261	1.248	1.0	122	0.00

Data Path : U:\HPCHEM1\BNA F\DATA\BF082117\
Data File : BF097909.D
Acq On : 21 Aug 2017 17:02
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 22 04:59:30 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 21 15:59:04 2017
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	1.184	1.173	0.9	120	0.00
89 C	Benzo(a)pyrene	1.116	1.097	1.7	119	0.00
90	Dibenzo(a,h)anthracene	0.909	0.813	10.6	108	0.00
91	Benzo(a,h,i)perylene	0.948	0.832	12.2	107	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 1