

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
 Data File : BG024070.D
 Acq On : 22 Sep 2016 4:40
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Sep 22 07:56:08 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 19:01:00 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	192#	0.00
2	1,4-Dioxane	0.470	0.476	-1.3	178#	0.00
3	Pyridine	1.412	1.440	-2.0	177#	0.00
4	n-Nitrosodimethylamine	0.744	0.755	-1.5	195#	0.00
5 S	2-Fluorophenol	1.139	1.227	-7.7	192#	0.00
6	Aniline	2.328	2.334	-0.3	186#	0.00
7 S	Phenol-d6	1.754	1.776	-1.3	184#	0.00
8	2-Chlorophenol	1.320	1.286	2.6	179#	0.00
9	Benzaldehyde	1.246	1.176	5.6	171#	0.00
10 C	Phenol	1.845	1.863	-1.0	181#	0.00
11	bis(2-Chloroethyl)ether	1.451	1.409	2.9	193#	0.00
12	1,3-Dichlorobenzene	1.545	1.537	0.5	191#	0.00
13 C	1,4-Dichlorobenzene	1.636	1.599	2.3	194#	0.00
14	1,2-Dichlorobenzene	1.536	1.495	2.7	188#	0.00
15	Benzyl Alcohol	1.546	1.494	3.4	172#	0.00
16	2,2'-oxybis(1-Chloropropane	1.945	1.897	2.5	193#	0.00
17	2-Methylphenol	1.243	1.180	5.1	174#	0.00
18	Hexachloroethane	0.630	0.631	-0.2	190#	0.00
19 P	n-Nitroso-di-n-propylamine	1.549	1.403	9.4	173#	0.00
20	3+4-Methylphenols	1.732	1.644	5.1	168#	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	155#	0.00
22	Acetophenone	0.518	0.547	-5.6	164#	0.00
23 S	Nitrobenzene-d5	0.448	0.509	-13.6	181#	0.00
24	Nitrobenzene	0.482	0.546	-13.3	182#	0.00
25	Isophorone	0.869	0.903	-3.9	164#	0.00
26 C	2-Nitrophenol	0.183	0.191	-4.4	165#	0.00
27	2,4-Dimethylphenol	0.311	0.322	-3.5	151#	0.00
28	bis(2-Chloroethoxy)methane	0.439	0.461	-5.0	168#	0.00
29 C	2,4-Dichlorophenol	0.330	0.351	-6.4	162#	0.00
30	1,2,4-Trichlorobenzene	0.362	0.391	-8.0	175#	0.00
31	Naphthalene	1.031	1.022	0.9	162#	0.00
32	Benzoic acid	0.253	0.214	15.4	130	0.00
33	4-Chloroaniline	0.430	0.421	2.1	149	0.00
34 C	Hexachlorobutadiene	0.272	0.306	-12.5	173#	0.00
35	Caprolactam	0.116	0.104	10.3	134	0.00
36 C	4-Chloro-3-methylphenol	0.441	0.432	2.0	150	0.00
37	2-Methylnaphthalene	0.784	0.732	6.6	144	0.00
38 I	Acenaphthene-d10	1.000	1.000	0.0	129	0.00
39	1,2,4,5-Tetrachlorobenzene	0.664	0.780	-17.5	150	0.00
40 P	Hexachlorocyclopentadiene	0.335	0.045#	86.6#	18#	0.00
41 S	2,4,6-Tribromophenol	0.360	0.312	13.3	106	0.00
42 C	2,4,6-Trichlorophenol	0.443	0.483	-9.0	137	0.00
43	2,4,5-Trichlorophenol	0.448	0.478	-6.7	131	0.00
44 S	2-Fluorobiphenyl	1.395	1.458	-4.5	137	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
 Data File : BG024070.D
 Acq On : 22 Sep 2016 4:40
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Sep 22 07:56:08 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 19:01:00 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	1.609	1.686	-4.8	136	0.00
46	2-Chloronaphthalene	1.181	1.253	-6.1	140	0.00
47	2-Nitroaniline	0.465	0.528	-13.5	149	0.00
48	Acenaphthylene	1.969	1.972	-0.2	129	0.00
49	Dimethylphthalate	1.715	1.615	5.8	123	0.00
50	2,6-Dinitrotoluene	0.362	0.351	3.0	125	0.00
51 C	Acenaphthene	1.217	1.203	1.2	130	0.00
52	3-Nitroaniline	0.339	0.355	-4.7	128	0.00
53 P	2,4-Dinitrophenol	0.226	0.032#	85.8#	17#	0.01
54	Dibenzofuran	1.900	1.845	2.9	126	0.00
55 P	4-Nitrophenol	0.269	0.235	12.6	117	0.00
56	2,4-Dinitrotoluene	0.532	0.493	7.3	118	0.00
57	Fluorene	1.647	1.538	6.6	122	0.00
58	2,3,4,6-Tetrachlorophenol	0.455	0.429	5.7	120	0.00
59	Diethylphthalate	1.831	1.638	10.5	116	0.00
60	4-Chlorophenyl-phenylether	0.887	0.819	7.7	117	0.00
61	4-Nitroaniline	0.342	0.344	-0.6	128	0.00
62	Azobenzene	1.773	1.737	2.0	128	0.00
63 I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
64	4,6-Dinitro-2-methylphenol	0.139	0.039	71.9#	31#	0.00
65 c	n-Nitrosodiphenylamine	0.568	0.588	-3.5	115	0.00
66	4-Bromophenyl-phenylether	0.243	0.246	-1.2	111	0.00
67	Hexachlorobenzene	0.285	0.273	4.2	105	0.00
68	Atrazine	0.244	0.251	-2.9	111	0.00
69 C	Pentachlorophenol	0.152	0.131	13.8	90	0.00
70	Phenanthrene	1.040	1.024	1.5	112	0.00
71	Anthracene	1.061	1.046	1.4	111	0.00
72	Carbazole	0.963	0.956	0.7	112	0.00
73	Di-n-butylphthalate	1.149	1.180	-2.7	117	0.00
74 C	Fluoranthene	1.253	1.235	1.4	109	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	110	0.00
76	Benzidine	0.619	0.626	-1.1	108	0.00
77	Pyrene	1.230	1.195	2.8	107	0.00
78 S	Terphenyl-d14	0.878	0.856	2.5	108	0.00
79	Butylbenzylphthalate	0.474	0.522	-10.1	128	0.00
80	Benzo(a)anthracene	1.120	1.155	-3.1	116	0.00
81	3,3'-Dichlorobenzidine	0.448	0.471	-5.1	114	0.00
82	Chrysene	1.066	1.085	-1.8	115	0.00
83	Bis(2-ethylhexyl)phthalate	0.649	0.707	-8.9	130	0.00
84 c	Di-n-octyl phthalate	1.065	1.198	-12.5	133	0.00
85	Indeno(1,2,3-cd)pyrene	1.180	0.935	20.8	92	0.03
86 I	Perylene-d12	1.000	1.000	0.0	110	0.00
87	Benzo(b)fluoranthene	1.136	1.184	-4.2	116	0.00

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092116\
Data File : BG024070.D
Acq On : 22 Sep 2016 4:40
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Sep 22 07:56:08 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 20 19:01:00 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	1.127	1.248	-10.7	122	0.00
89 C	Benzo(a)pyrene	1.079	1.162	-7.7	121	0.00
90	Dibenzo(a,h)anthracene	1.061	0.883	16.8	93	0.02
91	Benzo(a,h,i)perylene	1.020	0.682	33.1#	75	0.03

(#) = Out of Range

SPCC's out = 2 CCC's out = 0