

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG071320\
 Data File : BG046017.D
 Acq On : 13 Jul 2020 9:14
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 13 11:01:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:58:53 2020
 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	96	0.00
2	1,4-Dioxane	0.569	0.563	1.1	96	0.00
3	Pyridine	1.720	1.736	-0.9	98	0.00
4	n-Nitrosodimethylamine	0.582	0.601	-3.3	96	0.00
5 S	2-Fluorophenol	1.233	1.141	7.5	92	0.00
6	Aniline	2.271	2.232	1.7	96	0.00
7 S	Phenol-d6	1.776	1.730	2.6	95	0.00
8	2-Chlorophenol	1.328	1.232	7.2	92	0.00
9	Benzaldehyde	1.125	1.016	9.7	100	0.00
10 C	Phenol	1.781	1.729	2.9	96	0.00
11	bis(2-Chloroethyl)ether	1.473	1.426	3.2	96	0.00
12	1,3-Dichlorobenzene	1.530	1.470	3.9	95	0.00
13 C	1,4-Dichlorobenzene	1.505	1.468	2.5	95	0.00
14	1,2-Dichlorobenzene	1.449	1.391	4.0	94	0.00
15	Benzyl Alcohol	1.392	1.322	5.0	92	0.00
16	2,2'-oxybis(1-Chloropropane	1.922	1.868	2.8	94	0.00
17	2-Methylphenol	1.336	1.279	4.3	93	0.00
18	Hexachloroethane	0.579	0.538	7.1	91	0.00
19 P	n-Nitroso-di-n-propylamine	1.247	1.219	2.2	96	0.00
20	3+4-Methylphenols	1.821	1.772	2.7	93	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22	Acetophenone	0.534	0.511	4.3	95	0.00
23 S	Nitrobenzene-d5	0.347	0.364	-4.9	104	0.00
24	Nitrobenzene	0.384	0.400	-4.2	101	0.00
25	Isophorone	0.848	0.830	2.1	96	0.00
26 C	2-Nitrophenol	0.125	0.123	1.6	100	0.00
27	2,4-Dimethylphenol	0.302	0.283	6.3	92	0.00
28	bis(2-Chloroethoxy)methane	0.476	0.466	2.1	96	0.00
29 C	2,4-Dichlorophenol	0.313	0.291	7.0	91	0.00
30	1,2,4-Trichlorobenzene	0.346	0.332	4.0	95	0.00
31	Naphthalene	1.072	1.030	3.9	96	0.00
32	Benzoic acid	0.163	0.174	-6.7	101	0.00
33	4-Chloroaniline	0.479	0.457	4.6	94	0.00
34 C	Hexachlorobutadiene	0.202	0.194	4.0	96	0.00
35	Caprolactam	0.119	0.120	-0.8	98	0.00
36 C	4-Chloro-3-methylphenol	0.356	0.352	1.1	96	0.00
37	2-Methylnaphthalene	0.762	0.752	1.3	97	0.00
38	1-Methylnaphthalene	0.719	0.706	1.8	97	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
40	1,2,4,5-Tetrachlorobenzene	0.564	0.530	6.0	97	0.00
41 P	Hexachlorocyclopentadiene	0.335	0.279	16.7	85	0.00
42 S	2,4,6-Tribromophenol	0.186	0.181	2.7	96	0.00
43 C	2,4,6-Trichlorophenol	0.381	0.354	7.1	94	0.00
44	2,4,5-Trichlorophenol	0.431	0.416	3.5	97	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG071320\
 Data File : BG046017.D
 Acq On : 13 Jul 2020 9:14
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 13 11:01:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:58:53 2020
 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 S	2-Fluorobiphenyl	1.282	1.218	5.0	99	0.00
46	1,1'-Biphenyl	1.505	1.433	4.8	98	0.00
47	2-Chloronaphthalene	1.187	1.127	5.1	98	0.00
48	2-Nitroaniline	0.278	0.352	-26.6#	112	0.00
49	Acenaphthylene	1.915	1.842	3.8	98	0.00
50	Dimethylphthalate	1.501	1.460	2.7	100	0.00
51	2,6-Dinitrotoluene	0.233	0.274	-17.6	109	0.00
52 C	Acenaphthene	1.208	1.174	2.8	99	0.00
53	3-Nitroaniline	0.289	0.343	-18.7	108	0.00
54 P	2,4-Dinitrophenol	0.082	0.088	-7.3	101	0.00
55	Dibenzofuran	1.771	1.741	1.7	101	0.00
56 P	4-Nitrophenol	0.245	0.259	-5.7	106	0.00
57	2,4-Dinitrotoluene	0.288	0.364	-26.4#	113	0.00
58	Fluorene	1.453	1.437	1.1	101	0.00
59	2,3,4,6-Tetrachlorophenol	0.333	0.327	1.8	97	0.00
60	Diethylphthalate	1.574	1.521	3.4	99	0.00
61	4-Chlorophenyl-phenylether	0.723	0.707	2.2	101	0.00
62	4-Nitroaniline	0.310	0.368	-18.7	115	0.00
63	Azobenzene	1.689	1.664	1.5	101	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
65	4,6-Dinitro-2-methylphenol	0.059	0.060	-1.7	106	0.00
66 c	n-Nitrosodiphenylamine	0.620	0.581	6.3	100	0.00
67	4-Bromophenyl-phenylether	0.223	0.213	4.5	100	0.00
68	Hexachlorobenzene	0.224	0.211	5.8	102	0.00
69	Atrazine	0.183	0.165	9.8	94	0.00
70 C	Pentachlorophenol	0.125	0.117	6.4	93	0.00
71	Phenanthrene	1.057	1.014	4.1	102	0.00
72	Anthracene	1.073	1.019	5.0	102	0.00
73	Carbazole	1.076	1.036	3.7	103	0.00
74	Di-n-butylphthalate	1.300	1.198	7.8	101	0.00
75 C	Fluoranthene	1.222	1.182	3.3	105	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	104	0.00
77	Benzidine	0.572	0.618	-8.0	112	0.00
78	Pyrene	1.365	1.367	-0.1	104	0.00
79 S	Terphenyl-d14	0.899	0.871	3.1	104	0.00
80	Butylbenzylphthalate	0.645	0.633	1.9	100	0.00
81	Benzo(a)anthracene	1.315	1.251	4.9	101	0.00
82	3,3'-Dichlorobenzidine	0.486	0.452	7.0	98	0.00
83	Chrysene	1.253	1.194	4.7	100	0.00
84	Bis(2-ethylhexyl)phthalate	0.936	0.880	6.0	99	0.00
85 c	Di-n-octyl phthalate	1.625	1.487	8.5	96	0.00
86 I	Perylene-d12	1.000	1.000	0.0	103	0.00
87	Indeno(1,2,3-cd)pyrene	1.463	1.438	1.7	100	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG071320\
Data File : BG046017.D
Acq On : 13 Jul 2020 9:14
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jul 13 11:01:19 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 13 10:58:53 2020
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(b)fluoranthene	1.270	1.250	1.6	100	0.00
89	Benzo(k)fluoranthene	1.200	1.196	0.3	101	0.00
90 C	Benzo(a)pyrene	1.194	1.182	1.0	101	0.00
91	Dibenzo(a,h)anthracene	1.180	1.145	3.0	99	0.00
92	Benzo(q,h,i)perylene	1.188	1.152	3.0	99	0.00

(#) = Out of Range

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