Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016632.D
Acq On : 22 Apr 2015 22:34
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 23 03:21:12 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 23 01:45:23 2015
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	127	0.00
2		1,4-Dioxane	0.561	0.525	6.4	116	0.00
3		Pyridine	1.552	1.537	1.0	120	0.00
4		n-Nitrosodimethylamine	0.465	0.469	-0.9	125	0.00
5	S	2-Fluorophenol	1.248	1.254	-0.5	125	0.00
6		Aniline	2.408	2.551	-5.9	129	0.00
7	S	Phenol-d6	1.743	1.842	-5.7	128	0.00
8		2-Chlorophenol	1.490	1.566	-5.1	131	0.00
9		Benzaldehyde	0.793	0.739	6.8	133	0.00
10	C	Phenol	1.844	1.947	-5.6	127	0.00
11		bis(2-Chloroethyl)ether	1.444	1.494	-3.5	131	0.00
12		1,3-Dichlorobenzene	1.555	1.580	-1.6	130	0.00
13	C	1,4-Dichlorobenzene	1.583	1.613	-1.9	129	0.00
14		1,2-Dichlorobenzene	1.506	1.557	-3.4	131	0.00
15		Benzyl Alcohol	1.109	1.246	-12.4	133	0.00
16		2,2'-oxybis(1-Chloropropane	1.398	1.477	-5.7	136	0.00
17		2-Methylphenol	1.243	1.338	-7.6	132	0.00
18		Hexachloroethane	0.569	0.593	-4.2	133	0.00
19	P	n-Nitroso-di-n-propylamine	1.127	1.246	-10.6	137	0.00
20		3+4-Methylphenols	1.712	1.901	-11.0	132	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	135	0.00
22		Acetophenone	0.445	0.452	-1.6	132	0.00
23	S	Nitrobenzene-d5	0.325	0.330	-1.5	131	0.00
24		Nitrobenzene	0.338	0.344	-1.8	132	0.00
25		Isophorone	0.665	0.706	-6.2	134	0.00
26	C	2-Nitrophenol	0.187	0.206	-10.2	137	0.00
27		2,4-Dimethylphenol	0.317	0.334	-5.4	134	0.00
28		bis(2-Chloroethoxy)methane	0.399	0.414	-3.8	137	0.00
29	C	2,4-Dichlorophenol	0.262	0.291	-11.1	138	0.00
30		1,2,4-Trichlorobenzene	0.279	0.288	-3.2	136	0.00
31		Naphthalene	1.044	1.054	-1.0	132	0.00
32		Benzoic acid	0.103	0.171	-66.0#	208#	0.02
33		4-Chloroaniline	0.477	0.510	-6.9	132	0.00
34	C	Hexachlorobutadiene	0.124	0.129	-4.0	144	0.00
35		Caprolactam	0.152	0.161	-5.9	120	0.02
36	C	4-Chloro-3-methylphenol	0.317	0.344	-8.5	131	0.00
37		2-Methylnaphthalene	0.747	0.782	-4.7	135	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	128	0.00
39		1,2,4,5-Tetrachlorobenzene	0.460	0.486	-5.7	139	0.00
40	P	Hexachlorocyclopentadiene	0.143	0.155	-8.4	142	0.00
41	S	2,4,6-Tribromophenol	0.136	0.147	-8.1	121	0.00
42	C	2,4,6-Trichlorophenol	0.326	0.376	-15.3	139	0.00
43		2,4,5-Trichlorophenol	0.347	0.389	-12.1	133	0.00
44	S	2-Fluorobiphenyl	1.208	1.243	-2.9	133	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016632.D
Acq On : 22 Apr 2015 22:34
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 23 03:21:12 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 23 01:45:23 2015
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	1.540	1.604	-4.2	133	0.00
46		2-Chloronaphthalene	1.180	1.230	-4.2	132	0.00
47		2-Nitroaniline	0.344	0.362	-5.2	119	0.00
48		Acenaphthylene	2.094	2.145	-2.4	126	0.00
49		Dimethylphthalate	1.479	1.502	-1.6	122	0.00
50		2,6-Dinitrotoluene	0.363	0.379	-4.4	123	0.00
51	C	Acenaphthene	1.250	1.278	-2.2	128	0.00
52		3-Nitroaniline	0.449	0.462	-2.9	115	0.00
53	P	2,4-Dinitrophenol	0.133	0.121	9.0	99	0.00
54		Dibenzofuran	1.759	1.759	0.0	122	0.00
55	P	4-Nitrophenol	0.340	0.322	5.3	105	0.01
56		2,4-Dinitrotoluene	0.500	0.518	-3.6	113	0.00
57		Fluorene	1.550	1.564	-0.9	121	0.00
58		2,3,4,6-Tetrachlorophenol	0.262	0.288	-9.9	124	0.00
59		Diethylphthalate	1.576	1.556	1.3	115	0.00
60		4-Chlorophenyl-phenylether	0.636	0.640	-0.6	124	0.00
61		4-Nitroaniline	0.514	0.507	1.4	106	0.00
62		Azobenzene	1.462	1.455	0.5	117	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
64		4,6-Dinitro-2-methylphenol	0.129	0.134	-3.9	106	0.00
65	c	n-Nitrosodiphenylamine	0.671	0.719	-7.2	117	0.00
66		4-Bromophenyl-phenylether	0.157	0.173	-10.2	121	0.00
67		Hexachlorobenzene	0.164	0.175	-6.7	117	0.00
68		Atrazine	0.187	0.193	-3.2	107	0.00
69	C	Pentachlorophenol	0.073	0.085	-16.4	118	0.00
70		Phenanthrene	1.139	1.153	-1.2	109	0.00
71		Anthracene	1.131	1.160	-2.6	109	0.00
72		Carbazole	1.154	1.150	0.3	104	0.00
73		Di-n-butylphthalate	1.405	1.395	0.7	103	0.00
74	C	Fluoranthene	1.226	1.188	3.1	103	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76		Benzidine	0.740	0.738	0.3	100	0.00
77		Pyrene	1.343	1.435	-6.9	102	0.00
78	S	Terphenyl-d14	0.830	0.871	-4.9	104	0.00
79		Butylbenzylphthalate	0.703	0.761	-8.3	103	0.00
80		Benzo(a)anthracene	1.206	1.246	-3.3	103	0.00
81		3,3'-Dichlorobenzidine	0.399	0.422	-5.8	104	0.00
82		Chrysene	1.158	1.178	-1.7	101	0.00
83		Bis(2-ethylhexyl)phthalate	0.965	1.045	-8.3	103	0.00
84	c	Di-n-octyl phthalate	1.603	1.744	-8.8	103	0.00
85		Indeno(1,2,3-cd)pyrene	1.271	1.319	-3.8	103	0.00
86	I	Perylene-d12	1.000	1.000	0.0	103	0.00
87		Benzo(b)fluoranthene	1.193	1.205	-1.0	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042215\
Data File : BG016632.D
Acq On : 22 Apr 2015 22:34
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 23 03:21:12 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 23 01:45:23 2015
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	1.171	1.235	-5.5	103	0.00
89 C	Benzo(a)pyrene	1.133	1.176	-3.8	103	0.00
90	Dibenzo(a,h)anthracene	1.101	1.140	-3.5	102	0.00
91	Benzo(g,h,i)perylene	1.111	1.149	-3.4	101	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0