Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060515\
Data File : BG017473.D
Acq On : 5 Jun 2015 14:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 05 22:51:19 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 05 22:50:24 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	116	0.00
2		1,4-Dioxane	0.511	0.499	2.3	119	-0.02
3		Pyridine	1.365	1.275	6.6	104	-0.02
4		n-Nitrosodimethylamine	0.895	1.047	-17.0	130	-0.02
5	S	2-Fluorophenol	1.135	1.150	-1.3	117	0.00
6		Aniline	2.076	2.006	3.4	112	0.00
7	S	Phenol-d6	1.547	1.509	2.5	111	0.00
8		2-Chlorophenol	1.342	1.306	2.7	114	0.00
9		Benzaldehyde	1.050	1.004	4.4	106	-0.01
10	C	Phenol	1.589	1.602	-0.8	112	0.01
11		bis(2-Chloroethyl)ether	1.210	1.174	3.0	111	0.00
12		1,3-Dichlorobenzene	1.501	1.492	0.6	116	0.00
13	C	1,4-Dichlorobenzene	1.573	1.558	1.0	115	0.00
14		1,2-Dichlorobenzene	1.472	1.444	1.9	114	0.00
15		Benzyl Alcohol	1.440	1.468	-1.9	113	0.00
16		2,2'-oxybis(1-Chloropropane	2.040	1.903	6.7	107	0.00
17		2-Methylphenol	1.205	1.173	2.7	111	0.01
18		Hexachloroethane	0.627	0.624	0.5	113	0.00
19	P	n-Nitroso-di-n-propylamine	1.308	1.201	8.2	105	0.00
20		3+4-Methylphenols	1.668	1.657	0.7	112	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
22		Acetophenone	0.490	0.516	-5.3	115	0.00
23	S	Nitrobenzene-d5	0.401	0.423	-5.5	114	0.00
24		Nitrobenzene	0.416	0.438	-5.3	115	0.00
25		Isophorone	0.838	0.800	4.5	106	0.00
26	C	2-Nitrophenol	0.193	0.198	-2.6	109	0.00
27		2,4-Dimethylphenol	0.314	0.326	-3.8	114	0.00
28		bis(2-Chloroethoxy)methane	0.383	0.382	0.3	107	0.00
29	C	2,4-Dichlorophenol	0.306	0.333	-8.8	117	0.02
30		1,2,4-Trichlorobenzene	0.359	0.381	-6.1	118	0.00
31		Naphthalene	1.045	1.048	-0.3	110	0.00
32		Benzoic acid	0.198	0.214	-8.1	115	0.03
33		4-Chloroaniline	0.464	0.451	2.8	106	0.00
34	C	Hexachlorobutadiene	0.234	0.256	-9.4	120	0.00
35		Caprolactam	0.168	0.157	6.5	97	0.00
36	C	4-Chloro-3-methylphenol	0.398	0.386	3.0	103	0.02
37		2-Methylnaphthalene	0.803	0.811	-1.0	110	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
39		1,2,4,5-Tetrachlorobenzene	0.609	0.624	-2.5	116	0.00
40	P	Hexachlorocyclopentadiene	0.304	0.315	-3.6	109	0.00
41	S	2,4,6-Tribromophenol	0.278	0.284	-2.2	111	0.00
42	C	2,4,6-Trichlorophenol	0.436	0.470	-7.8	112	0.00
43		2,4,5-Trichlorophenol	0.477	0.498	-4.4	116	0.03
44	S	2-Fluorobiphenyl	1.326	1.366	-3.0	113	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060515\
Data File : BG017473.D
Acq On : 5 Jun 2015 14:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 05 22:51:19 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 05 22:50:24 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	1.462	1.491	-2.0	111	0.00
46		2-Chloronaphthalene	1.222	1.248	-2.1	108	0.00
47		2-Nitroaniline	0.451	0.449	0.4	106	0.00
48		Acenaphthylene	2.153	2.193	-1.9	110	0.00
49		Dimethylphthalate	1.741	1.674	3.8	104	0.00
50		2,6-Dinitrotoluene	0.391	0.378	3.3	105	0.00
51	C	Acenaphthene	1.269	1.300	-2.4	113	0.00
52		3-Nitroaniline	0.424	0.396	6.6	101	0.00
53	P	2,4-Dinitrophenol	0.205	0.207	-1.0	106	0.00
54		Dibenzofuran	1.873	1.856	0.9	107	0.00
55	P	4-Nitrophenol	0.306	0.302	1.3	97	0.05
56		2,4-Dinitrotoluene	0.557	0.563	-1.1	105	0.00
57		Fluorene	1.698	1.712	-0.8	109	0.00
58		2,3,4,6-Tetrachlorophenol	0.435	0.446	-2.5	106	0.01
59		Diethylphthalate	1.890	1.759	6.9	101	0.00
60		4-Chlorophenyl-phenylether	0.860	0.890	-3.5	113	0.00
61		4-Nitroaniline	0.466	0.483	-3.6	106	0.00
62		Azobenzene	1.796	1.776	1.1	106	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
64		4,6-Dinitro-2-methylphenol	0.148	0.150	-1.4	105	0.00
65	c	n-Nitrosodiphenylamine	0.597	0.598	-0.2	103	0.00
66		4-Bromophenyl-phenylether	0.221	0.229	-3.6	111	0.00
67		Hexachlorobenzene	0.238	0.250	-5.0	111	0.00
68		Atrazine	0.263	0.268	-1.9	104	0.01
69	C	Pentachlorophenol	0.128	0.135	-5.5	106	0.01
70		Phenanthrene	1.115	1.122	-0.6	107	0.00
71		Anthracene	1.114	1.126	-1.1	107	0.00
72		Carbazole	1.100	1.095	0.5	105	0.00
73		Di-n-butylphthalate	1.409	1.407	0.1	104	0.01
74	C	Fluoranthene	1.458	1.455	0.2	106	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	109	0.01
76		Benzidine	0.687	0.587	14.6	87	0.01
77		Pyrene	1.232	1.217	1.2	106	0.01
78	S	Terphenyl-d14	0.968	0.969	-0.1	108	0.00
79		Butylbenzylphthalate	0.533	0.543	-1.9	107	0.01
80		Benzo(a)anthracene	1.230	1.255	-2.0	109	0.00
81		3,3'-Dichlorobenzidine	0.460	0.471	-2.4	108	0.01
82		Chrysene	1.114	1.140	-2.3	110	0.00
83		Bis(2-ethylhexyl)phthalate	0.769	0.788	-2.5	109	0.01
84	c	Di-n-octyl phthalate	1.289	1.302	-1.0	109	0.01
85		Indeno(1,2,3-cd)pyrene	1.399	1.455	-4.0	111	0.04
86	I	Perylene-d12	1.000	1.000	0.0	109	0.02
87		Benzo(b)fluoranthene	1.267	1.292	-2.0	107	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060515\
Data File : BG017473.D
Acq On : 5 Jun 2015 14:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 05 22:51:19 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 05 22:50:24 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	1.213	1.251	-3.1	114	0.02
89 C	Benzo(a)pyrene	1.162	1.200	-3.3	111	0.02
90	Dibenzo(a,h)anthracene	1.208	1.260	-4.3	112	0.02
91	Benzo(g,h,i)perylene	1.150	1.186	-3.1	112	0.04

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