Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060615\
Data File : BG017521.D
Acq On : 7 Jun 2015 7:03
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 08 09:08:27 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 04 01:31:15 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	121	0.00
2		1,4-Dioxane	0.511	0.419	18.0	104	-0.02
3		Pyridine	1.365	1.273	6.7	108	-0.02
4		n-Nitrosodimethylamine	0.895	1.047	-17.0	135	-0.02
5	S	2-Fluorophenol	1.135	1.091	3.9	115	0.00
6		Aniline	2.076	2.009	3.2	116	0.00
7	S	Phenol-d6	1.547	1.502	2.9	114	0.00
8		2-Chlorophenol	1.342	1.304	2.8	118	0.00
9		Benzaldehyde	1.050	1.011	3.7	111	0.00
10	C	Phenol	1.589	1.578	0.7	115	0.00
11		bis(2-Chloroethyl)ether	1.210	1.226	-1.3	120	0.00
12		1,3-Dichlorobenzene	1.501	1.464	2.5	119	0.00
13	C	1,4-Dichlorobenzene	1.573	1.482	5.8	114	0.00
14		1,2-Dichlorobenzene	1.472	1.465	0.5	120	0.00
15		Benzyl Alcohol	1.440	1.422	1.3	114	0.00
16		2,2'-oxybis(1-Chloropropane	2.040	2.021	0.9	118	0.00
17		2-Methylphenol	1.205	1.200	0.4	118	0.00
18		Hexachloroethane	0.627	0.648	-3.3	122	0.00
19	P	n-Nitroso-di-n-propylamine	1.308	1.286	1.7	117	0.00
20		3+4-Methylphenols	1.668	1.645	1.4	115	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	118	0.00
22		Acetophenone	0.490	0.494	-0.8	116	0.00
23	S	Nitrobenzene-d5	0.401	0.428	-6.7	122	0.00
24		Nitrobenzene	0.416	0.439	-5.5	122	0.00
25		Isophorone	0.838	0.814	2.9	114	0.00
26	C	2-Nitrophenol	0.193	0.200	-3.6	117	0.00
27		2,4-Dimethylphenol	0.314	0.308	1.9	113	0.00
28		bis(2-Chloroethoxy)methane	0.383	0.399	-4.2	118	0.00
29	C	2,4-Dichlorophenol	0.306	0.319	-4.2	119	0.00
30		1,2,4-Trichlorobenzene	0.359	0.360	-0.3	118	0.00
31		Naphthalene	1.045	1.072	-2.6	119	0.00
32		Benzoic acid	0.198	0.194	2.0	110	0.03
33		4-Chloroaniline	0.464	0.458	1.3	114	0.00
34	C	Hexachlorobutadiene	0.234	0.254	-8.5	126	0.00
35		Caprolactam	0.168	0.145	13.7	94	0.00
36	C	4-Chloro-3-methylphenol	0.398	0.397	0.3	112	0.00
37		2-Methylnaphthalene	0.803	0.812	-1.1	116	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	119	0.00
39		1,2,4,5-Tetrachlorobenzene	0.609	0.626	-2.8	124	0.00
40	P	Hexachlorocyclopentadiene	0.304	0.316	-3.9	116	0.00
41	S	2,4,6-Tribromophenol	0.278	0.261	6.1	108	0.00
42	C	2,4,6-Trichlorophenol	0.436	0.429	1.6	109	0.00
43		2,4,5-Trichlorophenol	0.477	0.474	0.6	117	0.00
44	S	2-Fluorobiphenyl	1.326	1.401	-5.7	124	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060615\
Data File : BG017521.D
Acq On : 7 Jun 2015 7:03
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 08 09:08:27 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 04 01:31:15 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	1.528	-4.5	121	0.00
46		2-Chloronaphthalene	1.222	1.254	-2.6	116	0.00
47		2-Nitroaniline	0.451	0.443	1.8	111	0.00
48		Acenaphthylene	2.153	2.159	-0.3	115	0.00
49		Dimethylphthalate	1.741	1.642	5.7	108	0.00
50		2,6-Dinitrotoluene	0.391	0.361	7.7	107	0.00
51	C	Acenaphthene	1.269	1.275	-0.5	118	0.00
52		3-Nitroaniline	0.424	0.385	9.2	105	0.00
53	P	2,4-Dinitrophenol	0.205	0.145	29.3#	79	0.00
54		Dibenzofuran	1.873	1.843	1.6	113	0.00
55	P	4-Nitrophenol	0.306	0.277	9.5	94	0.02
56		2,4-Dinitrotoluene	0.557	0.522	6.3	104	0.00
57		Fluorene	1.698	1.712	-0.8	116	0.00
58		2,3,4,6-Tetrachlorophenol	0.435	0.433	0.5	110	0.00
59		Diethylphthalate	1.890	1.759	6.9	108	0.00
60		4-Chlorophenyl-phenylether	0.860	0.889	-3.4	120	0.00
61		4-Nitroaniline	0.466	0.394	15.5	92	0.00
62		Azobenzene	1.796	1.881	-4.7	120	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
64		4,6-Dinitro-2-methylphenol	0.148	0.134	9.5	95	0.00
65	c	n-Nitrosodiphenylamine	0.597	0.622	-4.2	109	0.00
66		4-Bromophenyl-phenylether	0.221	0.232	-5.0	114	0.00
67		Hexachlorobenzene	0.238	0.248	-4.2	112	0.00
68		Atrazine	0.263	0.259	1.5	102	0.00
69	C	Pentachlorophenol	0.128	0.118	7.8	94	0.00
70		Phenanthrene	1.115	1.121	-0.5	108	0.00
71		Anthracene	1.114	1.129	-1.3	109	0.00
72		Carbazole	1.100	1.042	5.3	102	0.00
73		Di-n-butylphthalate	1.409	1.377	2.3	103	0.00
74	C	Fluoranthene	1.458	1.432	1.8	106	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
76		Benzidine	0.687	0.686	0.1	99	0.00
77		Pyrene	1.232	1.216	1.3	103	0.00
78	S	Terphenyl-d14	0.968	0.992	-2.5	108	0.00
79		Butylbenzylphthalate	0.533	0.547	-2.6	106	0.00
80		Benzo(a)anthracene	1.230	1.243	-1.1	106	0.00
81		3,3'-Dichlorobenzidine	0.460	0.451	2.0	101	0.00
82		Chrysene	1.114	1.126	-1.1	106	0.00
83		Bis(2-ethylhexyl)phthalate	0.769	0.775	-0.8	105	0.00
84	c	Di-n-octyl phthalate	1.289	1.270	1.5	104	0.00
85		Indeno(1,2,3-cd)pyrene	1.399	1.379	1.4	103	0.01
86	I	Perylene-d12	1.000	1.000	0.0	103	0.00
87		Benzo(b)fluoranthene	1.267	1.287	-1.6	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060615\
Data File : BG017521.D
Acq On : 7 Jun 2015 7:03
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 08 09:08:27 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 04 01:31:15 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	1.238	-2.1	107	0.00
89 C	Benzo(a)pyrene	1.162	1.181	-1.6	103	0.00
90	Dibenzo(a,h)anthracene	1.208	1.237	-2.4	103	0.00
91	Benzo(g,h,i)perylene	1.150	1.166	-1.4	104	0.01

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