Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG050615\
Data File : BG016911.D
Acq On : 6 May 2015 21:57
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 07 04:23:24 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 07 03:57:31 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	107	0.00
2		1,4-Dioxane	0.481	0.458	4.8	108	0.00
3		Pyridine	1.488	1.400	5.9	103	0.00
4		n-Nitrosodimethylamine	0.888	0.866	2.5	112	0.00
5	S	2-Fluorophenol	1.149	1.109	3.5	108	0.00
6		Aniline	2.304	2.178	5.5	106	0.00
7	S	Phenol-d6	1.641	1.598	2.6	109	0.00
8		2-Chlorophenol	1.331	1.302	2.2	111	0.00
9		Benzaldehyde	1.144	1.120	2.1	106	0.00
10	C	Phenol	1.760	1.719	2.3	110	0.00
11		bis(2-Chloroethyl)ether	1.361	1.304	4.2	106	0.00
12		1,3-Dichlorobenzene	1.467	1.352	7.8	105	0.00
13	C	1,4-Dichlorobenzene	1.486	1.366	8.1	103	0.00
14		1,2-Dichlorobenzene	1.426	1.298	9.0	102	0.00
15		Benzyl Alcohol	1.499	1.433	4.4	107	0.00
16		2,2'-oxybis(1-Chloropropane	2.519	2.364	6.2	107	0.00
17		2-Methylphenol	1.239	1.153	6.9	104	0.00
18		Hexachloroethane	0.611	0.567	7.2	104	-0.01
19	P	n-Nitroso-di-n-propylamine	1.439	1.349	6.3	107	0.00
20		3+4-Methylphenols	1.708	1.630	4.6	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
22		Acetophenone	0.476	0.442	7.1	106	0.00
23	S	Nitrobenzene-d5	0.409	0.398	2.7	113	0.00
24		Nitrobenzene	0.413	0.400	3.1	111	0.00
25		Isophorone	0.827	0.754	8.8	106	0.00
26	C	2-Nitrophenol	0.168	0.164	2.4	112	0.00
27		2,4-Dimethylphenol	0.305	0.280	8.2	104	0.00
28		bis(2-Chloroethoxy)methane	0.417	0.377	9.6	107	0.00
29	C	2,4-Dichlorophenol	0.288	0.258	10.4	103	0.00
30		1,2,4-Trichlorobenzene	0.305	0.275	9.8	103	0.00
31		Naphthalene	0.995	0.898	9.7	105	0.00
32		Benzoic acid	0.179	0.179	0.0	117	0.00
33		4-Chloroaniline	0.459	0.426	7.2	105	0.00
34	C	Hexachlorobutadiene	0.191	0.167	12.6	102	0.00
35		Caprolactam	0.150	0.137	8.7	105	0.00
36	C	4-Chloro-3-methylphenol	0.369	0.341	7.6	107	0.00
37		2-Methylnaphthalene	0.758	0.658	13.2	102	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
39		1,2,4,5-Tetrachlorobenzene	0.541	0.503	7.0	102	0.00
40	P	Hexachlorocyclopentadiene	0.349	0.301	13.8	95	0.00
41	S	2,4,6-Tribromophenol	0.201	0.194	3.5	106	0.00
42	C	2,4,6-Trichlorophenol	0.383	0.368	3.9	100	0.00
43		2,4,5-Trichlorophenol	0.407	0.394	3.2	104	0.00
44	S	2-Fluorobiphenyl	1.325	1.235	6.8	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG050615\
Data File : BG016911.D
Acq On : 6 May 2015 21:57
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 07 04:23:24 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 07 03:57:31 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	1.432	1.335	6.8	102	0.00
46		2-Chloronaphthalene	1.134	1.074	5.3	103	0.00
47		2-Nitroaniline	0.395	0.433	-9.6	117	0.00
48		Acenaphthylene	1.975	1.878	4.9	103	-0.01
49		Dimethylphthalate	1.493	1.390	6.9	102	0.00
50		2,6-Dinitrotoluene	0.310	0.311	-0.3	105	0.00
51	C	Acenaphthene	1.175	1.092	7.1	102	0.00
52		3-Nitroaniline	0.353	0.371	-5.1	109	0.00
53	P	2,4-Dinitrophenol	0.142	0.132	7.0	107	0.00
54		Dibenzofuran	1.669	1.579	5.4	102	0.00
55	P	4-Nitrophenol	0.299	0.296	1.0	106	0.00
56		2,4-Dinitrotoluene	0.397	0.432	-8.8	115	0.00
57		Fluorene	1.478	1.369	7.4	101	0.00
58		2,3,4,6-Tetrachlorophenol	0.351	0.339	3.4	101	0.00
59		Diethylphthalate	1.577	1.478	6.3	103	0.00
60		4-Chlorophenyl-phenylether	0.742	0.675	9.0	101	0.00
61		4-Nitroaniline	0.409	0.406	0.7	107	0.00
62		Azobenzene	1.633	1.580	3.2	106	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
64		4,6-Dinitro-2-methylphenol	0.099	0.102	-3.0	113	0.00
65	c	n-Nitrosodiphenylamine	0.611	0.566	7.4	102	0.00
66		4-Bromophenyl-phenylether	0.201	0.184	8.5	101	0.00
67		Hexachlorobenzene	0.212	0.200	5.7	105	0.00
68		Atrazine	0.220	0.199	9.5	97	0.00
69	C	Pentachlorophenol	0.137	0.130	5.1	100	0.00
70		Phenanthrene	1.029	0.950	7.7	101	0.00
71		Anthracene	1.038	0.947	8.8	100	0.00
72		Carbazole	0.987	0.911	7.7	99	0.00
73		Di-n-butylphthalate	1.280	1.196	6.6	101	-0.01
74	C	Fluoranthene	1.194	1.073	10.1	96	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	97	0.00
76		Benzidine	0.680	0.578	15.0	83	0.00
77		Pyrene	1.180	1.113	5.7	98	0.00
78	S	Terphenyl-d14	0.853	0.823	3.5	98	0.00
79		Butylbenzylphthalate	0.561	0.546	2.7	100	-0.01
80		Benzo(a)anthracene	1.073	1.022	4.8	98	0.00
81		3,3'-Dichlorobenzidine	0.419	0.390	6.9	95	-0.01
82		Chrysene	1.005	0.952	5.3	97	0.00
83		Bis(2-ethylhexyl)phthalate	0.767	0.746	2.7	98	-0.01
84	c	Di-n-octyl phthalate	1.289	1.264	1.9	99	-0.01
85		Indeno(1,2,3-cd)pyrene	1.198	1.145	4.4	101	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	99	-0.01
87		Benzo(b)fluoranthene	1.114	1.042	6.5	98	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG050615\
Data File : BG016911.D
Acq On : 6 May 2015 21:57
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 07 04:23:24 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 07 03:57:31 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	1.072	1.006	6.2	99	-0.01
89 C	Benzo(a)pyrene	1.039	0.971	6.5	98	-0.01
90	Dibenzo(a,h)anthracene	1.061	0.999	5.8	101	-0.03
91	Benzo(g,h,i)perylene	1.011	0.947	6.3	102	-0.01

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