Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG050115\
Data File : BG016858.D
Acq On : 1 May 2015 22:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 02 00:06:35 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 01 23:47:41 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	109	0.00
2		1,4-Dioxane	0.465	0.448	3.7	109	0.00
3		Pyridine	1.305	1.309	-0.3	113	0.00
4		n-Nitrosodimethylamine	0.381	0.478	-25.5#	139	0.00
5	S	2-Fluorophenol	1.086	0.998	8.1	104	0.00
6		Aniline	2.060	2.074	-0.7	112	0.00
7	S	Phenol-d6	1.524	1.645	-7.9	120	0.00
8		2-Chlorophenol	1.380	1.408	-2.0	113	0.00
9		Benzaldehyde	0.869	0.906	-4.3	115	0.00
10	C	Phenol	1.563	1.562	0.1	111	0.00
11		bis(2-Chloroethyl)ether	1.247	1.252	-0.4	115	0.00
12		1,3-Dichlorobenzene	1.470	1.448	1.5	113	0.00
13	C	1,4-Dichlorobenzene	1.503	1.457	3.1	111	0.00
14		1,2-Dichlorobenzene	1.439	1.433	0.4	115	0.00
15		Benzyl Alcohol	0.950	0.861	9.4	100	0.00
16		2,2'-oxybis(1-Chloropropane	1.101	1.331	-20.9	137	0.00
17		2-Methylphenol	1.126	1.094	2.8	107	0.00
18		Hexachloroethane	0.530	0.583	-10.0	126	0.00
19	P	n-Nitroso-di-n-propylamine	0.974	1.085	-11.4	124	0.00
20		3+4-Methylphenols	1.547	1.640	-6.0	116	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.401	0.408	-1.7	118	0.00
23	S	Nitrobenzene-d5	0.307	0.328	-6.8	123	0.00
24		Nitrobenzene	0.290	0.311	-7.2	125	0.00
25		Isophorone	0.611	0.634	-3.8	122	0.00
26	C	2-Nitrophenol	0.184	0.176	4.3	110	0.00
27		2,4-Dimethylphenol	0.304	0.279	8.2	106	0.00
28		bis(2-Chloroethoxy)methane	0.356	0.358	-0.6	118	0.00
29	C	2,4-Dichlorophenol	0.264	0.247	6.4	108	0.00
30		1,2,4-Trichlorobenzene	0.280	0.265	5.4	112	0.00
31		Naphthalene	0.998	0.966	3.2	115	0.00
32		Benzoic acid	0.151	0.117	22.5	86	0.00
33		4-Chloroaniline	0.434	0.419	3.5	110	0.00
34	C	Hexachlorobutadiene	0.127	0.121	4.7	115	0.00
35		Caprolactam	0.145	0.126	13.1	100	0.00
36	C	4-Chloro-3-methylphenol	0.298	0.273	8.4	107	0.00
37		2-Methylnaphthalene	0.731	0.693	5.2	113	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
39		1,2,4,5-Tetrachlorobenzene	0.454	0.448	1.3	108	0.00
40	P	Hexachlorocyclopentadiene	0.080	0.127	-58.8#	169#	0.00
41	S	2,4,6-Tribromophenol	0.149	0.103	30.9#	75	0.00
42	C	2,4,6-Trichlorophenol	0.343	0.318	7.3	100	0.00
43		2,4,5-Trichlorophenol	0.348	0.340	2.3	103	0.00
44	S	2-Fluorobiphenyl	1.292	1.324	-2.5	114	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG050115\
Data File : BG016858.D
Acq On : 1 May 2015 22:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 02 00:06:35 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 01 23:47:41 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	1.469	1.459	0.7	110	0.00
46		2-Chloronaphthalene	1.134	1.135	-0.1	110	0.00
47		2-Nitroaniline	0.283	0.313	-10.6	120	0.00
48		Acenaphthylene	2.016	1.971	2.2	108	0.00
49		Dimethylphthalate	1.434	1.322	7.8	102	0.00
50		2,6-Dinitrotoluene	0.351	0.320	8.8	99	0.00
51	C	Acenaphthene	1.199	1.161	3.2	108	0.00
52		3-Nitroaniline	0.397	0.391	1.5	104	0.00
53	P	2,4-Dinitrophenol	0.123	0.110	10.6	96	0.00
54		Dibenzofuran	1.708	1.625	4.9	106	0.00
55	P	4-Nitrophenol	0.246	0.225	8.5	98	0.00
56		2,4-Dinitrotoluene	0.478	0.441	7.7	99	0.00
57		Fluorene	1.501	1.459	2.8	108	0.00
58		2,3,4,6-Tetrachlorophenol	0.265	0.254	4.2	102	0.00
59		Diethylphthalate	1.482	1.343	9.4	101	0.00
60		4-Chlorophenyl-phenylether	0.634	0.596	6.0	105	0.00
61		4-Nitroaniline	0.400	0.391	2.3	99	0.00
62		Azobenzene	1.226	1.023	16.6	93	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	68	0.00
64		4,6-Dinitro-2-methylphenol	0.118	0.166	-40.7#	94	0.00
65	c	n-Nitrosodiphenylamine	0.649	0.776	-19.6	83	0.00
66		4-Bromophenyl-phenylether	0.161	0.159	1.2	69	0.00
67		Hexachlorobenzene	0.168	0.168	0.0	70	0.00
68		Atrazine	0.189	0.174	7.9	64	0.00
69	C	Pentachlorophenol	0.059	0.067	-13.6	75	0.00
70		Phenanthrene	1.082	1.081	0.1	71	0.00
71		Anthracene	1.082	1.114	-3.0	72	0.00
72		Carbazole	1.058	1.064	-0.6	70	0.00
73		Di-n-butylphthalate	1.309	1.339	-2.3	72	0.00
74	C	Fluoranthene	1.163	1.113	4.3	68	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	70	0.00
76		Benzidine	0.769	0.616	19.9	54	0.00
77		Pyrene	1.308	1.251	4.4	69	0.00
78	S	Terphenyl-d14	0.829	0.874	-5.4	77	0.00
79		Butylbenzylphthalate	0.652	0.637	2.3	71	0.00
80		Benzo(a)anthracene	1.155	1.129	2.3	72	0.00
81		3,3'-Dichlorobenzidine	0.393	0.409	-4.1	74	0.00
82		Chrysene	1.093	1.064	2.7	72	0.00
83		Bis(2-ethylhexyl)phthalate	0.917	0.932	-1.6	73	0.00
84	c	Di-n-octyl phthalate	1.550	1.919	-23.8#	90	0.00
85		Indeno(1,2,3-cd)pyrene	1.244	1.640	-31.8#	99	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	94	0.00
87		Benzo(b)fluoranthene	1.139	1.084	4.8	94	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG050115\
Data File : BG016858.D
Acq On : 1 May 2015 22:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 02 00:06:35 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 01 23:47:41 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	1.115	1.096	1.7	95	0.03
89 C	Benzo(a)pyrene	1.078	1.052	2.4	96	0.00
90	Dibenzo(a,h)anthracene	1.088	1.060	2.6	99	0.00
91	Benzo(g,h,i)perylene	1.093	1.020	6.7	95	-0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 1