Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG022316\
Data File : BG021039.D
Acq On : 23 Feb 2016 23:07
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 24 04:04:07 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 24 00:35:06 2016
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	65	0.00
2		1,4-Dioxane	0.295	0.300	-1.7	66	0.00
3		Pyridine	1.043	0.994	4.7	64	0.00
4		n-Nitrosodimethylamine	0.429	0.491	-14.5	73	0.00
5	S	2-Fluorophenol	1.141	1.111	2.6	62	0.00
6		Aniline	2.072	2.120	-2.3	65	0.00
7	S	Phenol-d6	1.757	1.720	2.1	61	0.00
8		2-Chlorophenol	1.369	1.354	1.1	62	0.00
9		Benzaldehyde	0.850	0.886	-4.2	63	0.00
10	C	Phenol	1.833	1.841	-0.4	62	0.00
11		bis(2-Chloroethyl)ether	1.471	1.455	1.1	65	0.00
12		1,3-Dichlorobenzene	1.550	1.583	-2.1	66	0.00
13	C	1,4-Dichlorobenzene	1.618	1.609	0.6	65	0.00
14		1,2-Dichlorobenzene	1.503	1.465	2.5	63	0.00
15		Benzyl Alcohol	1.196	1.287	-7.6	67	0.00
16		2,2'-oxybis(1-Chloropropane	2.477	2.527	-2.0	66	0.00
17		2-Methylphenol	1.247	1.255	-0.6	64	0.00
18		Hexachloroethane	0.581	0.594	-2.2	66	0.00
19	P	n-Nitroso-di-n-propylamine	1.236	1.301	-5.3	66	0.00
20		3+4-Methylphenols	1.729	1.765	-2.1	64	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	67	0.00
22		Acetophenone	0.463	0.451	2.6	65	0.00
23	S	Nitrobenzene-d5	0.341	0.351	-2.9	66	0.00
24		Nitrobenzene	0.370	0.382	-3.2	67	0.00
25		Isophorone	0.702	0.714	-1.7	65	0.00
26	C	2-Nitrophenol	0.164	0.166	-1.2	63	0.00
27		2,4-Dimethylphenol	0.296	0.293	1.0	66	0.00
28		bis(2-Chloroethoxy)methane	0.380	0.368	3.2	64	0.00
29	C	2,4-Dichlorophenol	0.271	0.275	-1.5	66	0.00
30		1,2,4-Trichlorobenzene	0.290	0.293	-1.0	66	0.00
31		Naphthalene	1.046	1.037	0.9	65	0.00
32		Benzoic acid	0.224	0.241	-7.6	70	0.00
33		4-Chloroaniline	0.419	0.429	-2.4	65	0.00
34	C	Hexachlorobutadiene	0.161	0.168	-4.3	67	0.00
35		Caprolactam	0.106	0.111	-4.7	66	-0.03
36	C	4-Chloro-3-methylphenol	0.343	0.352	-2.6	65	0.00
37		2-Methylnaphthalene	0.761	0.773	-1.6	67	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	71	0.00
39		1,2,4,5-Tetrachlorobenzene	0.553	0.531	4.0	66	0.00
40	P	Hexachlorocyclopentadiene	0.313	0.327	-4.5	69	0.00
41	S	2,4,6-Tribromophenol	0.240	0.266	-10.8	74	0.00
42	C	2,4,6-Trichlorophenol	0.393	0.374	4.8	65	0.00
43		2,4,5-Trichlorophenol	0.403	0.402	0.2	67	0.00
44	S	2-Fluorobiphenyl	1.437	1.386	3.5	67	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG022316\
Data File : BG021039.D
Acq On : 23 Feb 2016 23:07
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 24 04:04:07 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 24 00:35:06 2016
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	1.735	1.686	2.8	67	0.00
46		2-Chloronaphthalene	1.214	1.181	2.7	68	0.00
47		2-Nitroaniline	0.366	0.397	-8.5	71	0.00
48		Acenaphthylene	2.219	2.198	0.9	68	0.00
49		Dimethylphthalate	1.631	1.623	0.5	68	0.00
50		2,6-Dinitrotoluene	0.334	0.361	-8.1	71	0.00
51	C	Acenaphthene	1.276	1.298	-1.7	70	0.00
52		3-Nitroaniline	0.384	0.426	-10.9	72	0.00
53	P	2,4-Dinitrophenol	0.180	0.211	-17.2	76	0.00
54		Dibenzofuran	1.812	1.847	-1.9	70	0.00
55	P	4-Nitrophenol	0.331	0.352	-6.3	74	0.00
56		2,4-Dinitrotoluene	0.480	0.504	-5.0	70	0.00
57		Fluorene	1.712	1.769	-3.3	71	0.00
58		2,3,4,6-Tetrachlorophenol	0.337	0.362	-7.4	72	0.00
59		Diethylphthalate	1.788	1.884	-5.4	72	0.00
60		4-Chlorophenyl-phenylether	0.789	0.808	-2.4	72	0.00
61		4-Nitroaniline	0.402	0.431	-7.2	72	0.00
62		Azobenzene	1.550	1.674	-8.0	73	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	74	0.00
64		4,6-Dinitro-2-methylphenol	0.121	0.128	-5.8	73	0.00
65	c	n-Nitrosodiphenylamine	0.602	0.601	0.2	71	0.00
66		4-Bromophenyl-phenylether	0.210	0.216	-2.9	74	0.00
67		Hexachlorobenzene	0.230	0.239	-3.9	74	0.00
68		Atrazine	0.189	0.200	-5.8	76	0.00
69	C	Pentachlorophenol	0.133	0.146	-9.8	77	0.00
70		Phenanthrene	1.131	1.172	-3.6	75	0.00
71		Anthracene	1.127	1.171	-3.9	75	0.00
72		Carbazole	1.016	1.074	-5.7	76	0.00
73		Di-n-butylphthalate	1.238	1.365	-10.3	78	0.00
74	C	Fluoranthene	1.138	1.249	-9.8	78	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
76		Benzidine	1.017	0.863	15.1	82	0.00
77		Pyrene	2.285	1.842	19.4	82	0.00
78	S	Terphenyl-d14	1.541	1.292	16.2	83	0.00
79		Butylbenzylphthalate	0.806	0.768	4.7	93	0.00
80		Benzo(a)anthracene	1.199	1.236	-3.1	105	0.00
81		3,3'-Dichlorobenzidine	0.443	0.466	-5.2	107	0.00
82		Chrysene	1.119	1.160	-3.7	107	0.00
83		Bis(2-ethylhexyl)phthalate	0.880	0.904	-2.7	106	0.00
84	c	Di-n-octyl phthalate	1.252	1.321	-5.5	112	0.00
85		Indeno(1,2,3-cd)pyrene	1.006	1.101	-9.4	115	0.00
86	I	Perylene-d12	1.000	1.000	0.0	112	0.00
87		Benzo(b)fluoranthene	1.319	1.347	-2.1	112	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG022316\
Data File : BG021039.D
Acq On : 23 Feb 2016 23:07
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 24 04:04:07 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 24 00:35:06 2016
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	1.244	1.255	-0.9	110	0.00
89 C	Benzo(a)pyrene	1.181	1.249	-5.8	112	0.00
90	Dibenzo(a,h)anthracene	1.064	1.109	-4.2	113	-0.01
91	Benzo(g,h,i)perylene	1.062	1.127	-6.1	116	-0.03

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