Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
Data File : BG024126.D
Acq On : 24 Sep 2016 16:29
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 26 00:51:54 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG092316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 23 15:35:28 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	108	0.00
2		1,4-Dioxane	0.442	0.534	-20.8	117	0.00
3		Pyridine	1.562	1.782	-14.1	115	0.00
4		n-Nitrosodimethylamine	0.792	0.793	-0.1	106	0.00
5	S	2-Fluorophenol	1.154	1.267	-9.8	109	0.00
6		Aniline	2.728	2.731	-0.1	108	0.00
7	S	Phenol-d6	1.942	2.033	-4.7	111	0.00
8		2-Chlorophenol	1.382	1.450	-4.9	113	0.00
9		Benzaldehyde	1.058	1.132	-7.0	114	0.00
10	C	Phenol	2.042	2.127	-4.2	111	0.00
11		bis(2-Chloroethyl)ether	1.610	1.617	-0.4	106	0.00
12		1,3-Dichlorobenzene	1.541	1.625	-5.5	113	0.00
13	C	1,4-Dichlorobenzene	1.573	1.620	-3.0	110	0.00
14		1,2-Dichlorobenzene	1.531	1.580	-3.2	109	0.00
15		Benzyl Alcohol	1.880	1.802	4.1	105	0.00
16		2,2'-oxybis(1-Chloropropane	2.223	2.336	-5.1	112	0.00
17		2-Methylphenol	1.358	1.462	-7.7	115	0.00
18		Hexachloroethane	0.663	0.686	-3.5	109	0.00
19	P	n-Nitroso-di-n-propylamine	1.908	1.851	3.0	105	0.00
20		3+4-Methylphenols	1.972	2.033	-3.1	113	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	114	0.00
22		Acetophenone	0.577	0.563	2.4	108	0.00
23	S	Nitrobenzene-d5	0.507	0.498	1.8	109	0.00
24		Nitrobenzene	0.544	0.535	1.7	109	0.00
25		Isophorone	1.035	0.992	4.2	109	0.00
26	C	2-Nitrophenol	0.193	0.197	-2.1	112	0.00
27		2,4-Dimethylphenol	0.323	0.330	-2.2	113	0.00
28		bis(2-Chloroethoxy)methane	0.503	0.503	0.0	112	0.00
29	C	2,4-Dichlorophenol	0.336	0.347	-3.3	110	0.00
30		1,2,4-Trichlorobenzene	0.359	0.353	1.7	108	0.00
31		Naphthalene	1.013	1.020	-0.7	113	0.00
32		Benzoic acid	0.241	0.243	-0.8	109	0.00
33		4-Chloroaniline	0.452	0.460	-1.8	113	0.00
34	C	Hexachlorobutadiene	0.283	0.261	7.8	103	0.00
35		Caprolactam	0.137	0.123	10.2	106	0.00
36	C	4-Chloro-3-methylphenol	0.472	0.471	0.2	108	0.00
37		2-Methylnaphthalene	0.776	0.769	0.9	110	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	112	0.00
39		1,2,4,5-Tetrachlorobenzene	0.584	0.581	0.5	106	0.00
40	P	Hexachlorocyclopentadiene	0.220	0.223	-1.4	100	0.00
41	S	2,4,6-Tribromophenol	0.287	0.251	12.5	97	0.00
42	C	2,4,6-Trichlorophenol	0.378	0.386	-2.1	107	0.00
43		2,4,5-Trichlorophenol	0.389	0.392	-0.8	106	0.00
44	S	2-Fluorobiphenyl	1.190	1.189	0.1	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
Data File : BG024126.D
Acq On : 24 Sep 2016 16:29
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 26 00:51:54 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG092316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 23 15:35:28 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	1.368	1.411	-3.1	111	0.00
46		2-Chloronaphthalene	1.021	1.033	-1.2	110	0.00
47		2-Nitroaniline	0.475	0.468	1.5	106	0.00
48		Acenaphthylene	1.734	1.742	-0.5	111	0.00
49		Dimethylphthalate	1.534	1.478	3.7	106	0.00
50		2,6-Dinitrotoluene	0.322	0.318	1.2	106	0.00
51	C	Acenaphthene	1.051	1.041	1.0	107	0.00
52		3-Nitroaniline	0.322	0.322	0.0	111	0.00
53	P	2,4-Dinitrophenol	0.190	0.170	10.5	96	0.00
54		Dibenzofuran	1.637	1.613	1.5	108	0.00
55	P	4-Nitrophenol	0.225	0.197	12.4	96	0.00
56		2,4-Dinitrotoluene	0.490	0.468	4.5	106	0.00
57		Fluorene	1.425	1.374	3.6	107	0.00
58		2,3,4,6-Tetrachlorophenol	0.375	0.349	6.9	98	0.00
59		Diethylphthalate	1.652	1.517	8.2	105	0.00
60		4-Chlorophenyl-phenylether	0.763	0.710	6.9	104	0.00
61		4-Nitroaniline	0.333	0.303	9.0	103	0.00
62		Azobenzene	1.679	1.589	5.4	105	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	0.139	0.138	0.7	94	0.00
65	c	n-Nitrosodiphenylamine	0.547	0.613	-12.1	107	0.00
66		4-Bromophenyl-phenylether	0.232	0.238	-2.6	99	0.00
67		Hexachlorobenzene	0.261	0.264	-1.1	98	0.00
68		Atrazine	0.242	0.241	0.4	97	0.00
69	C	Pentachlorophenol	0.124	0.115	7.3	90	0.00
70		Phenanthrene	1.019	1.061	-4.1	102	0.00
71		Anthracene	1.046	1.070	-2.3	101	0.00
72		Carbazole	0.974	0.966	0.8	99	0.00
73		Di-n-butylphthalate	1.275	1.199	6.0	95	0.00
74	C	Fluoranthene	1.330	1.220	8.3	95	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	124	0.00
76		Benzidine	0.566	0.539	4.8	107	0.00
77		Pyrene	1.397	1.179	15.6	96	0.00
78	S	Terphenyl-d14	0.977	0.805	17.6	96	0.00
79		Butylbenzylphthalate	0.514	0.515	-0.2	120	0.00
80		Benzo(a)anthracene	1.139	1.158	-1.7	122	0.00
81		3,3'-Dichlorobenzidine	0.408	0.453	-11.0	131	0.00
82		Chrysene	1.070	1.100	-2.8	125	0.00
83		Bis(2-ethylhexyl)phthalate	0.614	0.735	-19.7	138	0.00
84	c	Di-n-octyl phthalate	1.060	1.279	-20.7#	135	0.00
85		Indeno(1,2,3-cd)pyrene	1.193	1.382	-15.8	130	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	135	-0.02
87		Benzo(b)fluoranthene	1.164	1.162	0.2	129	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
Data File : BG024126.D
Acq On : 24 Sep 2016 16:29
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 26 00:51:54 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG092316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 23 15:35:28 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	1.156	1.149	0.6	129	0.00
89 C	Benzo(a)pyrene	1.125	1.125	0.0	131	0.00
90	Dibenzo(a,h)anthracene	1.123	1.105	1.6	130	-0.03
91	Benzo(g,h,i)perylene	1.132	1.113	1.7	130	-0.02

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