Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030216\
Data File : BG021224.D
Acq On : 2 Mar 2016 22:19
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 03 00:14:10 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 11:46:16 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	100	-0.02
2		1,4-Dioxane	0.565	0.444	21.4	83	0.00
3		Pyridine	1.786	1.539	13.8	81	0.00
4		n-Nitrosodimethylamine	0.900	0.698	22.4	73	0.00
5	S	2-Fluorophenol	1.258	1.176	6.5	92	0.00
6		Aniline	2.742	2.370	13.6	84	0.00
7	S	Phenol-d6	2.069	1.874	9.4	87	0.00
8		2-Chlorophenol	1.542	1.483	3.8	93	0.00
9		Benzaldehyde	1.292	1.061	17.9	86	0.00
10	C	Phenol	2.205	1.986	9.9	88	0.00
11		bis(2-Chloroethyl)ether	1.700	1.511	11.1	84	0.00
12		1,3-Dichlorobenzene	1.581	1.558	1.5	97	-0.02
13	C	1,4-Dichlorobenzene	1.628	1.574	3.3	96	-0.02
14		1,2-Dichlorobenzene	1.553	1.530	1.5	95	0.00
15		Benzyl Alcohol	1.889	1.585	16.1	80	0.00
16		2,2'-oxybis(1-Chloropropane	2.934	2.346	20.0	79	0.00
17		2-Methylphenol	1.526	1.360	10.9	87	0.00
18		Hexachloroethane	0.682	0.644	5.6	93	-0.02
19	P	n-Nitroso-di-n-propylamine	1.911	1.488	22.1	76	-0.02
20		3+4-Methylphenols	2.161	1.899	12.1	86	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	90	0.00
22		Acetophenone	0.595	0.553	7.1	82	0.00
23	S	Nitrobenzene-d5	0.473	0.456	3.6	84	0.00
24		Nitrobenzene	0.516	0.495	4.1	85	0.00
25		Isophorone	1.024	0.867	15.3	75	0.00
26	C	2-Nitrophenol	0.188	0.193	-2.7	90	-0.02
27		2,4-Dimethylphenol	0.337	0.315	6.5	82	0.00
28		bis(2-Chloroethoxy)methane	0.476	0.431	9.5	81	0.00
29	C	2,4-Dichlorophenol	0.304	0.303	0.3	89	0.00
30		1,2,4-Trichlorobenzene	0.306	0.325	-6.2	94	0.00
31		Naphthalene	1.043	1.034	0.9	89	0.00
32		Benzoic acid	0.261	0.223	14.6	72	0.00
33		4-Chloroaniline	0.481	0.437	9.1	80	0.00
34	C	Hexachlorobutadiene	0.186	0.214	-15.1	101	-0.02
35		Caprolactam	0.148	0.112	24.3	67	0.01
36	C	4-Chloro-3-methylphenol	0.461	0.398	13.7	78	0.00
37		2-Methylnaphthalene	0.742	0.698	5.9	84	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	79	0.00
39		1,2,4,5-Tetrachlorobenzene	0.601	0.701	-16.6	89	0.00
40	P	Hexachlorocyclopentadiene	0.329	0.216	34.3#	48#	0.00
41	S	2,4,6-Tribromophenol	0.208	0.208	0.0	78	0.00
42	C	2,4,6-Trichlorophenol	0.451	0.480	-6.4	84	0.00
43		2,4,5-Trichlorophenol	0.482	0.499	-3.5	81	0.00
44	S	2-Fluorobiphenyl	1.408	1.518	-7.8	83	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030216\
Data File : BG021224.D
Acq On : 2 Mar 2016 22:19
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 03 00:14:10 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 11:46:16 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	1.645	1.770	-7.6	84	0.00
46		2-Chloronaphthalene	1.235	1.317	-6.6	82	0.00
47		2-Nitroaniline	0.602	0.546	9.3	71	0.00
48		Acenaphthylene	2.090	2.112	-1.1	78	0.00
49		Dimethylphthalate	1.782	1.664	6.6	74	0.00
50		2,6-Dinitrotoluene	0.373	0.360	3.5	75	0.00
51	C	Acenaphthene	1.327	1.260	5.0	72	0.00
52		3-Nitroaniline	0.414	0.370	10.6	70	0.00
53	P	2,4-Dinitrophenol	0.173	0.115	33.5#	45#	0.00
54		Dibenzofuran	1.772	1.730	2.4	77	0.00
55	P	4-Nitrophenol	0.343	0.317	7.6	71	0.00
56		2,4-Dinitrotoluene	0.537	0.501	6.7	72	0.00
57		Fluorene	1.689	1.632	3.4	75	0.00
58		2,3,4,6-Tetrachlorophenol	0.386	0.370	4.1	75	0.00
59		Diethylphthalate	1.970	1.765	10.4	71	0.00
60		4-Chlorophenyl-phenylether	0.843	0.827	1.9	77	0.00
61		4-Nitroaniline	0.442	0.401	9.3	70	0.00
62		Azobenzene	2.116	1.852	12.5	69	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	74	0.00
64		4,6-Dinitro-2-methylphenol	0.128	0.098	23.4	52	0.00
65	c	n-Nitrosodiphenylamine	0.615	0.636	-3.4	74	0.00
66		4-Bromophenyl-phenylether	0.209	0.221	-5.7	75	0.00
67		Hexachlorobenzene	0.203	0.222	-9.4	78	0.00
68		Atrazine	0.211	0.225	-6.6	77	0.00
69	C	Pentachlorophenol	0.133	0.135	-1.5	70	0.00
70		Phenanthrene	1.079	1.092	-1.2	73	0.00
71		Anthracene	1.104	1.116	-1.1	73	0.00
72		Carbazole	0.979	0.992	-1.3	73	0.00
73		Di-n-butylphthalate	1.377	1.354	1.7	70	0.00
74	C	Fluoranthene	1.268	1.275	-0.6	73	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	72	0.00
76		Benzidine	0.560	0.601	-7.3	75	0.00
77		Pyrene	1.193	1.213	-1.7	73	0.00
78	S	Terphenyl-d14	0.826	0.863	-4.5	74	0.00
79		Butylbenzylphthalate	0.609	0.584	4.1	67	0.00
80		Benzo(a)anthracene	1.184	1.212	-2.4	73	0.00
81		3,3'-Dichlorobenzidine	0.448	0.471	-5.1	74	0.00
82		Chrysene	1.122	1.148	-2.3	73	0.00
83		Bis(2-ethylhexyl)phthalate	0.892	0.867	2.8	69	0.00
84	c	Di-n-octyl phthalate	1.486	1.456	2.0	69	0.00
85		Indeno(1,2,3-cd)pyrene	1.287	1.231	4.4	71	0.00
86	I	Perylene-d12	1.000	1.000	0.0	76	0.00
87		Benzo(b)fluoranthene	1.258	1.267	-0.7	74	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030216\
Data File : BG021224.D
Acq On : 2 Mar 2016 22:19
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 03 00:14:10 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 02 11:46:16 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	1.260	1.244	1.3	74	0.00
89 C	Benzo(a)pyrene	1.216	1.230	-1.2	76	0.00
90	Dibenzo(a,h)anthracene	1.203	1.079	10.3	68	0.00
91	Benzo(g,h,i)perylene	1.162	0.972	16.4	65	0.00

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