Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019130.D
Acq On : 10 Oct 2015 23:08
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 12 06:40:07 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 12 06:30:50 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	95	0.00
2		1,4-Dioxane	0.413	0.386	6.5	97	0.00
3		Pyridine	1.263	1.235	2.2	95	0.00
4		n-Nitrosodimethylamine	0.716	0.688	3.9	93	0.00
5	S	2-Fluorophenol	0.982	0.932	5.1	94	0.00
6		Aniline	2.043	1.992	2.5	96	0.00
7	S	Phenol-d6	1.509	1.503	0.4	97	0.00
8		2-Chlorophenol	1.233	1.215	1.5	98	0.00
9		Benzaldehyde	0.950	0.939	1.2	97	0.00
10	C	Phenol	1.571	1.546	1.6	95	0.00
11		bis(2-Chloroethyl)ether	1.187	1.152	2.9	98	0.00
12		1,3-Dichlorobenzene	1.346	1.294	3.9	95	0.00
13	C	1,4-Dichlorobenzene	1.377	1.331	3.3	97	0.00
14		1,2-Dichlorobenzene	1.332	1.254	5.9	94	0.00
15		Benzyl Alcohol	1.348	1.320	2.1	93	0.00
16		2,2'-oxybis(1-Chloropropane	2.543	2.440	4.1	96	0.00
17		2-Methylphenol	1.122	1.108	1.2	94	0.00
18		Hexachloroethane	0.525	0.501	4.6	94	0.00
19	P	n-Nitroso-di-n-propylamine	1.197	1.163	2.8	94	0.00
20		3+4-Methylphenols	1.616	1.635	-1.2	97	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
22		Acetophenone	0.468	0.452	3.4	96	0.00
23	S	Nitrobenzene-d5	0.362	0.353	2.5	94	0.00
24		Nitrobenzene	0.384	0.373	2.9	95	0.00
25		Isophorone	0.736	0.710	3.5	93	0.00
26	C	2-Nitrophenol	0.164	0.161	1.8	95	0.00
27		2,4-Dimethylphenol	0.286	0.277	3.1	95	0.00
28		bis(2-Chloroethoxy)methane	0.393	0.372	5.3	93	0.00
29	C	2,4-Dichlorophenol	0.300	0.290	3.3	94	0.00
30		1,2,4-Trichlorobenzene	0.322	0.307	4.7	95	0.00
31		Naphthalene	0.963	0.922	4.3	95	0.00
32		Benzoic acid	0.158	0.140	11.4	78	0.00
33		4-Chloroaniline	0.447	0.417	6.7	92	0.00
34	C	Hexachlorobutadiene	0.221	0.217	1.8	96	0.00
35		Caprolactam	0.110	0.111	-0.9	94	0.00
36	C	4-Chloro-3-methylphenol	0.381	0.365	4.2	93	0.00
37		2-Methylnaphthalene	0.748	0.711	4.9	94	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
39		1,2,4,5-Tetrachlorobenzene	0.575	0.560	2.6	93	0.00
40	P	Hexachlorocyclopentadiene	0.299	0.300	-0.3	91	0.00
41	S	2,4,6-Tribromophenol	0.273	0.261	4.4	93	0.00
42	C	2,4,6-Trichlorophenol	0.400	0.396	1.0	91	0.00
43		2,4,5-Trichlorophenol	0.425	0.415	2.4	90	0.00
44	S	2-Fluorobiphenyl	1.301	1.251	3.8	94	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019130.D
Acq On : 10 Oct 2015 23:08
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 12 06:40:07 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 12 06:30:50 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	1.452	1.387	4.5	94	0.00
46		2-Chloronaphthalene	1.118	1.081	3.3	93	0.00
47		2-Nitroaniline	0.449	0.439	2.2	93	0.00
48		Acenaphthylene	1.971	1.875	4.9	92	0.00
49		Dimethylphthalate	1.580	1.472	6.8	92	0.00
50		2,6-Dinitrotoluene	0.337	0.326	3.3	92	0.00
51	C	Acenaphthene	1.185	1.093	7.8	90	0.00
52		3-Nitroaniline	0.372	0.343	7.8	90	0.00
53	P	2,4-Dinitrophenol	0.159	0.151	5.0	85	0.00
54		Dibenzofuran	1.751	1.643	6.2	92	0.00
55	P	4-Nitrophenol	0.282	0.273	3.2	90	0.00
56		2,4-Dinitrotoluene	0.491	0.476	3.1	93	0.00
57		Fluorene	1.519	1.436	5.5	94	0.00
58		2,3,4,6-Tetrachlorophenol	0.403	0.390	3.2	91	0.00
59		Diethylphthalate	1.641	1.508	8.1	90	0.00
60		4-Chlorophenyl-phenylether	0.774	0.712	8.0	91	0.00
61		4-Nitroaniline	0.405	0.384	5.2	91	0.00
62		Azobenzene	1.522	1.452	4.6	93	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
64		4,6-Dinitro-2-methylphenol	0.109	0.107	1.8	88	0.00
65	c	n-Nitrosodiphenylamine	0.553	0.533	3.6	92	0.00
66		4-Bromophenyl-phenylether	0.203	0.203	0.0	94	0.00
67		Hexachlorobenzene	0.239	0.235	1.7	93	0.00
68		Atrazine	0.228	0.220	3.5	91	0.00
69	C	Pentachlorophenol	0.124	0.125	-0.8	87	0.00
70		Phenanthrene	1.029	0.968	5.9	91	0.00
71		Anthracene	1.030	0.980	4.9	91	0.00
72		Carbazole	0.964	0.920	4.6	92	0.00
73		Di-n-butylphthalate	1.180	1.114	5.6	90	0.00
74	C	Fluoranthene	1.324	1.235	6.7	91	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	92	0.00
76		Benzidine	0.513	0.492	4.1	87	0.00
77		Pyrene	1.120	1.076	3.9	90	0.00
78	S	Terphenyl-d14	0.757	0.728	3.8	91	0.00
79		Butylbenzylphthalate	0.456	0.429	5.9	88	0.00
80		Benzo(a)anthracene	1.072	1.025	4.4	90	0.00
81		3,3'-Dichlorobenzidine	0.397	0.384	3.3	91	0.00
82		Chrysene	1.018	0.969	4.8	91	0.00
83		Bis(2-ethylhexyl)phthalate	0.634	0.598	5.7	89	0.00
84	c	Di-n-octyl phthalate	1.095	1.055	3.7	92	0.00
85		Indeno(1,2,3-cd)pyrene	1.238	1.208	2.4	93	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	95	-0.01
87		Benzo(b)fluoranthene	1.075	1.013	5.8	93	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019130.D
Acq On : 10 Oct 2015 23:08
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 12 06:40:07 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 12 06:30:50 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	1.062	1.009	5.0	93	-0.01
89 C	Benzo(a)pyrene	1.016	0.965	5.0	93	-0.01
90	Dibenzo(a,h)anthracene	1.091	1.033	5.3	93	-0.02
91	Benzo(g,h,i)perylene	1.014	0.956	5.7	93	-0.02

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