Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG121117\
Data File : BG031468.D
Acq On : 11 Dec 2017 23:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 12 04:41:52 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 11 17:25:55 2017
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	117	0.00
2		1,4-Dioxane	0.538	0.493	8.4	109	0.00
3		Pyridine	1.417	1.381	2.5	108	0.00
4		n-Nitrosodimethylamine	0.668	0.633	5.2	105	0.00
5	S	2-Fluorophenol	1.128	1.137	-0.8	112	0.00
6		Aniline	2.063	1.976	4.2	109	0.00
7	S	Phenol-d6	1.566	1.529	2.4	110	0.00
8		2-Chlorophenol	1.259	1.237	1.7	111	0.00
9		Benzaldehyde	0.908	0.844	7.0	106	0.00
10	C	Phenol	1.609	1.595	0.9	110	0.00
11		bis(2-Chloroethyl)ether	1.313	1.269	3.4	108	0.00
12		1,3-Dichlorobenzene	1.497	1.433	4.3	110	0.00
13	C	1,4-Dichlorobenzene	1.556	1.464	5.9	109	0.00
14		1,2-Dichlorobenzene	1.482	1.407	5.1	111	0.00
15		Benzyl Alcohol	1.225	1.104	9.9	103	0.00
16		2,2'-oxybis(1-Chloropropane	1.475	1.610	-9.2	126	0.00
17		2-Methylphenol	1.097	1.078	1.7	110	0.00
18		Hexachloroethane	0.524	0.505	3.6	110	0.00
19	P	n-Nitroso-di-n-propylamine	1.233	1.111	9.9	103	0.00
20		3+4-Methylphenols	1.634	1.503	8.0	106	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	114	0.00
22		Acetophenone	0.480	0.462	3.7	105	0.00
23	S	Nitrobenzene-d5	0.324	0.305	5.9	102	0.00
24		Nitrobenzene	0.333	0.314	5.7	102	0.00
25		Isophorone	0.613	0.584	4.7	101	0.00
26	C	2-Nitrophenol	0.164	0.160	2.4	106	0.00
27		2,4-Dimethylphenol	0.250	0.244	2.4	105	0.00
28		bis(2-Chloroethoxy)methane	0.396	0.387	2.3	107	0.00
29	C	2,4-Dichlorophenol	0.294	0.295	-0.3	106	0.00
30		1,2,4-Trichlorobenzene	0.376	0.356	5.3	108	0.00
31		Naphthalene	0.983	0.916	6.8	105	0.00
32		Benzoic acid	0.132	0.110	16.7	95	0.00
33		4-Chloroaniline	0.427	0.407	4.7	105	0.00
34	C	Hexachlorobutadiene	0.249	0.235	5.6	110	0.00
35		Caprolactam	0.116	0.113	2.6	107	0.00
36	C	4-Chloro-3-methylphenol	0.336	0.317	5.7	105	0.00
37		2-Methylnaphthalene	0.761	0.714	6.2	105	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	115	0.00
39		1,2,4,5-Tetrachlorobenzene	0.782	0.730	6.6	105	0.00
40	P	Hexachlorocyclopentadiene	0.182	0.155	14.8	94	0.00
41	S	2,4,6-Tribromophenol	0.234	0.237	-1.3	110	0.00
42	C	2,4,6-Trichlorophenol	0.399	0.393	1.5	105	0.00
43		2,4,5-Trichlorophenol	0.430	0.431	-0.2	107	0.00
44	S	2-Fluorobiphenyl	1.363	1.264	7.3	105	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG121117\
Data File : BG031468.D
Acq On : 11 Dec 2017 23:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 12 04:41:52 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 11 17:25:55 2017
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	1.643	1.530	6.9	105	0.00
46		2-Chloronaphthalene	1.202	1.114	7.3	105	0.00
47		2-Nitroaniline	0.338	0.311	8.0	101	0.00
48		Acenaphthylene	1.935	1.812	6.4	105	0.00
49		Dimethylphthalate	1.656	1.553	6.2	105	0.00
50		2,6-Dinitrotoluene	0.354	0.335	5.4	104	0.00
51	C	Acenaphthene	1.223	1.144	6.5	106	0.00
52		3-Nitroaniline	0.361	0.349	3.3	108	0.00
53	P	2,4-Dinitrophenol	0.116	0.099	14.7	90	0.00
54		Dibenzofuran	1.952	1.801	7.7	105	0.00
55	P	4-Nitrophenol	0.222	0.218	1.8	106	0.00
56		2,4-Dinitrotoluene	0.503	0.478	5.0	105	0.00
57		Fluorene	1.564	1.458	6.8	105	0.00
58		2,3,4,6-Tetrachlorophenol	0.404	0.400	1.0	105	0.00
59		Diethylphthalate	1.660	1.601	3.6	107	0.00
60		4-Chlorophenyl-phenylether	0.957	0.886	7.4	104	0.00
61		4-Nitroaniline	0.379	0.368	2.9	108	0.00
62		Azobenzene	1.375	1.278	7.1	103	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	113	0.00
64		4,6-Dinitro-2-methylphenol	0.112	0.107	4.5	103	0.00
65	c	n-Nitrosodiphenylamine	0.585	0.568	2.9	107	0.00
66		4-Bromophenyl-phenylether	0.233	0.226	3.0	108	0.00
67		Hexachlorobenzene	0.240	0.228	5.0	108	0.00
68		Atrazine	0.214	0.217	-1.4	110	0.00
69	C	Pentachlorophenol	0.100	0.096	4.0	102	0.00
70		Phenanthrene	1.045	0.985	5.7	107	0.00
71		Anthracene	1.033	0.978	5.3	106	0.00
72		Carbazole	0.935	0.910	2.7	108	0.00
73		Di-n-butylphthalate	1.080	1.082	-0.2	110	0.00
74	C	Fluoranthene	1.307	1.262	3.4	109	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	117	0.00
76		Benzidine	0.465	0.462	0.6	107	0.00
77		Pyrene	1.243	1.187	4.5	108	0.00
78	S	Terphenyl-d14	0.879	0.826	6.0	108	0.00
79		Butylbenzylphthalate	0.436	0.452	-3.7	112	0.00
80		Benzo(a)anthracene	1.207	1.148	4.9	109	0.00
81		3,3'-Dichlorobenzidine	0.420	0.434	-3.3	112	0.00
82		Chrysene	1.157	1.103	4.7	109	0.00
83		Bis(2-ethylhexyl)phthalate	0.627	0.638	-1.8	112	0.00
84	c	Di-n-octyl phthalate	1.035	1.066	-3.0	112	-0.01
85		Indeno(1,2,3-cd)pyrene	1.237	1.243	-0.5	113	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	119	-0.01
87		Benzo(b)fluoranthene	1.196	1.135	5.1	111	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG121117\
Data File : BG031468.D
Acq On : 11 Dec 2017 23:37
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 12 04:41:52 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 11 17:25:55 2017
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	1.220	1.156	5.2	111	-0.01
89 C	Benzo(a)pyrene	1.134	1.087	4.1	111	-0.01
90	Dibenzo(a,h)anthracene	1.085	1.057	2.6	113	-0.03
91	Benzo(g,h,i)perylene	1.066	1.032	3.2	112	-0.03

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