Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG092518\
Data File : BG036965.D
Acq On : 26 Sep 2018 2:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 26 02:37:42 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 24 10:41:23 2018
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	134	0.00
2		1,4-Dioxane	0.477	0.549	-15.1	144	0.00
3		Pyridine	1.378	1.546	-12.2	142	0.00
4		n-Nitrosodimethylamine	0.612	0.718	-17.3	153#	0.00
5	S	2-Fluorophenol	1.043	1.180	-13.1	147	0.00
6		Aniline	2.094	2.113	-0.9	132	0.00
7	S	Phenol-d6	1.613	1.705	-5.7	137	-0.01
8		2-Chlorophenol	1.211	1.307	-7.9	139	0.00
9		Benzaldehyde	1.066	1.051	1.4	129	0.00
10	C	Phenol	1.665	1.790	-7.5	139	0.00
11		bis(2-Chloroethyl)ether	1.540	1.477	4.1	132	-0.01
12		1,3-Dichlorobenzene	1.485	1.504	-1.3	132	-0.01
13	C	1,4-Dichlorobenzene	1.519	1.512	0.5	132	-0.01
14		1,2-Dichlorobenzene	1.472	1.456	1.1	133	0.00
15		Benzyl Alcohol	1.309	1.370	-4.7	137	0.00
16		2,2'-oxybis(1-Chloropropane	2.957	2.961	-0.1	131	-0.01
17		2-Methylphenol	1.160	1.172	-1.0	134	0.00
18		Hexachloroethane	0.559	0.583	-4.3	136	0.00
19	P	n-Nitroso-di-n-propylamine	1.342	1.290	3.9	127	0.00
20		3+4-Methylphenols	1.614	1.664	-3.1	134	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
22		Acetophenone	0.470	0.487	-3.6	132	0.00
23	S	Nitrobenzene-d5	0.373	0.396	-6.2	131	0.00
24		Nitrobenzene	0.399	0.412	-3.3	129	0.00
25		Isophorone	0.682	0.697	-2.2	128	0.00
26	C	2-Nitrophenol	0.159	0.182	-14.5	138	0.00
27		2,4-Dimethylphenol	0.248	0.258	-4.0	128	0.00
28		bis(2-Chloroethoxy)methane	0.421	0.418	0.7	123	0.00
29	C	2,4-Dichlorophenol	0.287	0.316	-10.1	133	-0.01
30		1,2,4-Trichlorobenzene	0.359	0.358	0.3	124	0.00
31		Naphthalene	0.952	0.937	1.6	125	-0.01
32		Benzoic acid	0.175	0.163	6.9	116	0.00
33		4-Chloroaniline	0.433	0.419	3.2	121	-0.01
34	C	Hexachlorobutadiene	0.248	0.252	-1.6	127	-0.01
35		Caprolactam	0.118	0.117	0.8	121	0.00
36	C	4-Chloro-3-methylphenol	0.343	0.368	-7.3	128	0.00
37		2-Methylnaphthalene	0.725	0.700	3.4	123	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	119	0.00
39		1,2,4,5-Tetrachlorobenzene	0.698	0.720	-3.2	123	0.00
40	P	Hexachlorocyclopentadiene	0.359	0.345	3.9	109	-0.01
41	S	2,4,6-Tribromophenol	0.239	0.249	-4.2	121	0.00
42	C	2,4,6-Trichlorophenol	0.387	0.438	-13.2	131	-0.01
43		2,4,5-Trichlorophenol	0.413	0.478	-15.7	131	0.00
44	S	2-Fluorobiphenyl	1.343	1.350	-0.5	119	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG092518\
Data File : BG036965.D
Acq On : 26 Sep 2018 2:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 26 02:37:42 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 24 10:41:23 2018
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	1.532	1.556	-1.6	122	0.00
46		2-Chloronaphthalene	1.205	1.222	-1.4	122	0.00
47		2-Nitroaniline	0.417	0.458	-9.8	126	0.00
48		Acenaphthylene	1.888	1.909	-1.1	121	0.00
49		Dimethylphthalate	1.610	1.580	1.9	117	0.00
50		2,6-Dinitrotoluene	0.351	0.354	-0.9	119	0.00
51	C	Acenaphthene	1.141	1.143	-0.2	120	0.00
52		3-Nitroaniline	0.370	0.385	-4.1	121	0.00
53	P	2,4-Dinitrophenol	0.136	0.145	-6.6	123	0.00
54		Dibenzofuran	1.930	1.917	0.7	118	0.00
55	P	4-Nitrophenol	0.236	0.261	-10.6	126	-0.01
56		2,4-Dinitrotoluene	0.490	0.495	-1.0	118	0.00
57		Fluorene	1.442	1.429	0.9	119	-0.01
58		2,3,4,6-Tetrachlorophenol	0.419	0.465	-11.0	126	0.00
59		Diethylphthalate	1.624	1.577	2.9	116	0.00
60		4-Chlorophenyl-phenylether	0.913	0.905	0.9	117	0.00
61		4-Nitroaniline	0.389	0.393	-1.0	117	0.00
62		Azobenzene	1.560	1.599	-2.5	122	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	117	0.00
64		4,6-Dinitro-2-methylphenol	0.105	0.109	-3.8	121	0.00
65	c	n-Nitrosodiphenylamine	0.552	0.565	-2.4	118	0.00
66		4-Bromophenyl-phenylether	0.227	0.230	-1.3	116	0.00
67		Hexachlorobenzene	0.234	0.230	1.7	113	0.00
68		Atrazine	0.224	0.218	2.7	110	-0.01
69	C	Pentachlorophenol	0.148	0.154	-4.1	116	-0.01
70		Phenanthrene	0.964	0.951	1.3	115	-0.01
71		Anthracene	0.953	0.954	-0.1	116	0.00
72		Carbazole	0.929	0.934	-0.5	116	0.00
73		Di-n-butylphthalate	1.032	1.054	-2.1	118	0.00
74	C	Fluoranthene	1.160	1.146	1.2	115	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	120	0.00
76		Benzidine	0.605	0.681	-12.6	136	0.00
77		Pyrene	1.200	1.149	4.2	114	-0.01
78	S	Terphenyl-d14	0.761	0.716	5.9	114	0.00
79		Butylbenzylphthalate	0.478	0.487	-1.9	123	-0.01
80		Benzo(a)anthracene	1.167	1.136	2.7	119	-0.01
81		3,3'-Dichlorobenzidine	0.444	0.444	0.0	117	0.00
82		Chrysene	1.128	1.094	3.0	118	0.00
83		Bis(2-ethylhexyl)phthalate	0.668	0.669	-0.1	123	-0.01
84	c	Di-n-octyl phthalate	1.100	1.120	-1.8	122	-0.02
85		Indeno(1,2,3-cd)pyrene	1.211	1.239	-2.3	120	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	121	-0.02
87		Benzo(b)fluoranthene	1.066	1.033	3.1	114	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG092518\
Data File : BG036965.D
Acq On : 26 Sep 2018 2:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 26 02:37:42 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 24 10:41:23 2018
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	1.069	1.037	3.0	119	-0.01
89 C	Benzo(a)pyrene	1.030	1.024	0.6	119	-0.01
90	Dibenzo(a,h)anthracene	0.966	0.974	-0.8	120	-0.02
91	Benzo(g,h,i)perylene	0.969	0.982	-1.3	119	-0.02

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