Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG092418\
Data File : BG036936.D
Acq On : 25 Sep 2018 3:35
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 25 07:29:03 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	120	0.00
2		1,4-Dioxane	0.477	0.530	-11.1	125	0.00
3		Pyridine	1.378	1.542	-11.9	128	0.00
4		n-Nitrosodimethylamine	0.612	0.677	-10.6	130	0.00
5	S	2-Fluorophenol	1.043	1.155	-10.7	129	0.00
6		Aniline	2.094	2.160	-3.2	121	0.00
7	S	Phenol-d6	1.613	1.813	-12.4	130	0.00
8		2-Chlorophenol	1.211	1.337	-10.4	128	0.00
9		Benzaldehyde	1.066	1.125	-5.5	124	0.00
10	C	Phenol	1.665	1.849	-11.1	129	0.00
11		bis(2-Chloroethyl)ether	1.540	1.541	-0.1	124	0.00
12		1,3-Dichlorobenzene	1.485	1.496	-0.7	118	0.00
13	C	1,4-Dichlorobenzene	1.519	1.531	-0.8	120	0.00
14		1,2-Dichlorobenzene	1.472	1.480	-0.5	122	0.00
15		Benzyl Alcohol	1.309	1.366	-4.4	123	0.00
16		2,2'-oxybis(1-Chloropropane	2.957	3.093	-4.6	123	-0.01
17		2-Methylphenol	1.160	1.236	-6.6	127	0.00
18		Hexachloroethane	0.559	0.588	-5.2	124	0.00
19	P	n-Nitroso-di-n-propylamine	1.342	1.396	-4.0	123	0.00
20		3+4-Methylphenols	1.614	1.758	-8.9	127	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	121	0.00
22		Acetophenone	0.470	0.465	1.1	121	0.00
23	S	Nitrobenzene-d5	0.373	0.388	-4.0	124	0.00
24		Nitrobenzene	0.399	0.405	-1.5	122	0.00
25		Isophorone	0.682	0.710	-4.1	126	0.00
26	C	2-Nitrophenol	0.159	0.172	-8.2	126	0.00
27		2,4-Dimethylphenol	0.248	0.247	0.4	118	0.00
28		bis(2-Chloroethoxy)methane	0.421	0.428	-1.7	121	0.00
29	C	2,4-Dichlorophenol	0.287	0.305	-6.3	123	0.00
30		1,2,4-Trichlorobenzene	0.359	0.360	-0.3	120	0.00
31		Naphthalene	0.952	0.946	0.6	121	-0.01
32		Benzoic acid	0.175	0.105	40.0#	72	-0.01
33		4-Chloroaniline	0.433	0.423	2.3	118	0.00
34	C	Hexachlorobutadiene	0.248	0.246	0.8	119	0.00
35		Caprolactam	0.118	0.118	0.0	118	0.00
36	C	4-Chloro-3-methylphenol	0.343	0.366	-6.7	122	0.00
37		2-Methylnaphthalene	0.725	0.713	1.7	120	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	121	0.00
39		1,2,4,5-Tetrachlorobenzene	0.698	0.693	0.7	119	0.00
40	P	Hexachlorocyclopentadiene	0.359	0.333	7.2	107	-0.01
41	S	2,4,6-Tribromophenol	0.239	0.246	-2.9	121	0.00
42	C	2,4,6-Trichlorophenol	0.387	0.418	-8.0	126	0.00
43		2,4,5-Trichlorophenol	0.413	0.452	-9.4	126	0.00
44	S	2-Fluorobiphenyl	1.343	1.320	1.7	118	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG092418\
Data File : BG036936.D
Acq On : 25 Sep 2018 3:35
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 25 07:29:03 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	1.515	1.1	120	0.00
46		2-Chloronaphthalene	1.205	1.218	-1.1	123	0.00
47		2-Nitroaniline	0.417	0.438	-5.0	122	0.00
48		Acenaphthylene	1.888	1.893	-0.3	121	0.00
49		Dimethylphthalate	1.610	1.565	2.8	117	0.00
50		2,6-Dinitrotoluene	0.351	0.351	0.0	120	0.00
51	C	Acenaphthene	1.141	1.157	-1.4	123	0.00
52		3-Nitroaniline	0.370	0.369	0.3	117	0.00
53	P	2,4-Dinitrophenol	0.136	0.053	61.0#	45#	0.00
54		Dibenzofuran	1.930	1.899	1.6	118	0.00
55	P	4-Nitrophenol	0.236	0.229	3.0	112	0.00
56		2,4-Dinitrotoluene	0.490	0.490	0.0	119	0.00
57		Fluorene	1.442	1.416	1.8	119	0.00
58		2,3,4,6-Tetrachlorophenol	0.419	0.447	-6.7	123	0.00
59		Diethylphthalate	1.624	1.564	3.7	117	0.00
60		4-Chlorophenyl-phenylether	0.913	0.894	2.1	117	0.00
61		4-Nitroaniline	0.389	0.383	1.5	116	0.00
62		Azobenzene	1.560	1.595	-2.2	123	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	118	0.00
64		4,6-Dinitro-2-methylphenol	0.105	0.076	27.6#	85	0.00
65	c	n-Nitrosodiphenylamine	0.552	0.563	-2.0	118	0.00
66		4-Bromophenyl-phenylether	0.227	0.230	-1.3	117	0.00
67		Hexachlorobenzene	0.234	0.235	-0.4	116	0.00
68		Atrazine	0.224	0.221	1.3	113	-0.01
69	C	Pentachlorophenol	0.148	0.140	5.4	107	0.00
70		Phenanthrene	0.964	0.960	0.4	117	0.00
71		Anthracene	0.953	0.965	-1.3	119	0.00
72		Carbazole	0.929	0.933	-0.4	117	0.00
73		Di-n-butylphthalate	1.032	1.038	-0.6	117	0.00
74	C	Fluoranthene	1.160	1.148	1.0	116	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	116	0.00
76		Benzidine	0.605	0.550	9.1	107	0.00
77		Pyrene	1.200	1.206	-0.5	116	0.00
78	S	Terphenyl-d14	0.761	0.736	3.3	113	0.00
79		Butylbenzylphthalate	0.478	0.482	-0.8	118	0.00
80		Benzo(a)anthracene	1.167	1.141	2.2	116	-0.01
81		3,3'-Dichlorobenzidine	0.444	0.439	1.1	112	0.00
82		Chrysene	1.128	1.100	2.5	115	0.00
83		Bis(2-ethylhexyl)phthalate	0.668	0.669	-0.1	119	-0.01
84	c	Di-n-octyl phthalate	1.100	1.099	0.1	116	-0.02
85		Indeno(1,2,3-cd)pyrene	1.211	1.254	-3.6	118	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	118	-0.02
87		Benzo(b)fluoranthene	1.066	1.065	0.1	116	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG092418\
Data File : BG036936.D
Acq On : 25 Sep 2018 3:35
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 25 07:29:03 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	1.050	1.8	118	-0.01
89 C	Benzo(a)pyrene	1.030	1.025	0.5	117	-0.01
90	Dibenzo(a,h)anthracene	0.966	0.986	-2.1	120	-0.02
91	Benzo(g,h,i)perylene	0.969	0.990	-2.2	118	-0.02

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