Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG091918\
Data File : BG036839.D
Acq On : 20 Sep 2018 7:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 20 09:01:42 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 31 13:03:20 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	124	-0.02
2		1,4-Dioxane	0.588	0.504	14.3	111	0.00
3		Pyridine	1.652	1.526	7.6	112	0.00
4		n-Nitrosodimethylamine	0.736	0.669	9.1	112	0.00
5	S	2-Fluorophenol	1.261	1.217	3.5	119	-0.01
6		Aniline	2.291	2.080	9.2	113	-0.01
7	S	Phenol-d6	1.805	1.742	3.5	119	-0.01
8		2-Chlorophenol	1.367	1.276	6.7	115	-0.01
9		Benzaldehyde	1.243	1.139	8.4	121	-0.01
10	C	Phenol	1.889	1.806	4.4	118	-0.01
11		bis(2-Chloroethyl)ether	1.498	1.385	7.5	116	-0.02
12		1,3-Dichlorobenzene	1.561	1.469	5.9	119	-0.02
13	C	1,4-Dichlorobenzene	1.576	1.485	5.8	118	-0.02
14		1,2-Dichlorobenzene	1.505	1.403	6.8	118	-0.02
15		Benzyl Alcohol	1.455	1.371	5.8	115	-0.01
16		2,2'-oxybis(1-Chloropropane	3.020	2.666	11.7	111	-0.01
17		2-Methylphenol	1.254	1.184	5.6	118	-0.02
18		Hexachloroethane	0.565	0.551	2.5	120	-0.02
19	P	n-Nitroso-di-n-propylamine	1.349	1.212	10.2	113	-0.01
20		3+4-Methylphenols	1.751	1.689	3.5	120	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	125	-0.02
22		Acetophenone	0.525	0.471	10.3	112	-0.01
23	S	Nitrobenzene-d5	0.369	0.387	-4.9	128	-0.01
24		Nitrobenzene	0.397	0.394	0.8	120	-0.01
25		Isophorone	0.732	0.665	9.2	113	-0.01
26	C	2-Nitrophenol	0.158	0.183	-15.8	144	-0.01
27		2,4-Dimethylphenol	0.292	0.264	9.6	114	-0.01
28		bis(2-Chloroethoxy)methane	0.449	0.415	7.6	115	-0.02
29	C	2,4-Dichlorophenol	0.320	0.325	-1.6	124	-0.01
30		1,2,4-Trichlorobenzene	0.374	0.361	3.5	123	-0.02
31		Naphthalene	1.000	0.929	7.1	116	-0.02
32		Benzoic acid	0.202	0.169	16.3	102	0.00
33		4-Chloroaniline	0.458	0.433	5.5	117	-0.01
34	C	Hexachlorobutadiene	0.251	0.251	0.0	123	-0.02
35		Caprolactam	0.127	0.120	5.5	113	0.00
36	C	4-Chloro-3-methylphenol	0.371	0.371	0.0	122	-0.01
37		2-Methylnaphthalene	0.732	0.694	5.2	118	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	127	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.708	0.683	3.5	123	-0.01
40	P	Hexachlorocyclopentadiene	0.368	0.349	5.2	119	-0.02
41	S	2,4,6-Tribromophenol	0.239	0.257	-7.5	129	-0.01
42	C	2,4,6-Trichlorophenol	0.431	0.449	-4.2	130	-0.01
43		2,4,5-Trichlorophenol	0.460	0.474	-3.0	127	-0.01
44	S	2-Fluorobiphenyl	1.401	1.326	5.4	122	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG091918\
Data File : BG036839.D
Acq On : 20 Sep 2018 7:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 20 09:01:42 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 31 13:03:20 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	1.660	1.517	8.6	119	-0.02
46		2-Chloronaphthalene	1.267	1.189	6.2	120	-0.01
47		2-Nitroaniline	0.398	0.436	-9.5	132	-0.01
48		Acenaphthylene	1.981	1.843	7.0	119	-0.02
49		Dimethylphthalate	1.633	1.543	5.5	120	-0.02
50		2,6-Dinitrotoluene	0.336	0.354	-5.4	132	-0.01
51	C	Acenaphthene	1.269	1.118	11.9	112	-0.01
52		3-Nitroaniline	0.362	0.383	-5.8	125	-0.01
53	P	2,4-Dinitrophenol	0.127	0.122	3.9	120	0.00
54		Dibenzofuran	1.987	1.860	6.4	120	-0.02
55	P	4-Nitrophenol	0.296	0.253	14.5	103	0.00
56		2,4-Dinitrotoluene	0.438	0.497	-13.5	140	-0.01
57		Fluorene	1.486	1.386	6.7	118	-0.01
58		2,3,4,6-Tetrachlorophenol	0.432	0.442	-2.3	126	-0.01
59		Diethylphthalate	1.646	1.534	6.8	117	-0.02
60		4-Chlorophenyl-phenylether	0.917	0.900	1.9	126	-0.02
61		4-Nitroaniline	0.379	0.388	-2.4	121	-0.01
62		Azobenzene	1.675	1.519	9.3	115	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	126	-0.02
64		4,6-Dinitro-2-methylphenol	0.088	0.097	-10.2	139	0.00
65	c	n-Nitrosodiphenylamine	0.594	0.548	7.7	119	-0.01
66		4-Bromophenyl-phenylether	0.234	0.230	1.7	125	-0.02
67		Hexachlorobenzene	0.239	0.237	0.8	125	0.00
68		Atrazine	0.223	0.208	6.7	121	-0.02
69	C	Pentachlorophenol	0.163	0.181	-11.0	131	0.00
70		Phenanthrene	1.016	0.932	8.3	117	-0.01
71		Anthracene	1.013	0.926	8.6	116	-0.01
72		Carbazole	1.012	0.896	11.5	111	-0.01
73		Di-n-butylphthalate	1.127	1.000	11.3	110	-0.02
74	C	Fluoranthene	1.239	1.107	10.7	111	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	116	-0.01
76		Benzidine	0.638	0.537	15.8	105	-0.01
77		Pyrene	1.230	1.173	4.6	110	0.00
78	S	Terphenyl-d14	0.856	0.824	3.7	113	-0.01
79		Butylbenzylphthalate	0.489	0.473	3.3	109	-0.01
80		Benzo(a)anthracene	1.212	1.120	7.6	109	-0.02
81		3,3'-Dichlorobenzidine	0.483	0.450	6.8	106	-0.01
82		Chrysene	1.148	1.074	6.4	109	-0.01
83		Bis(2-ethylhexyl)phthalate	0.685	0.637	7.0	106	-0.02
84	c	Di-n-octyl phthalate	1.144	1.058	7.5	104	-0.03
85		Indeno(1,2,3-cd)pyrene	1.334	1.188	10.9	103	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	114	-0.02
87		Benzo(b)fluoranthene	1.111	1.064	4.2	108	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG091918\
Data File : BG036839.D
Acq On : 20 Sep 2018 7:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 20 09:01:42 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 31 13:03:20 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	1.085	1.016	6.4	107	-0.02
89 C	Benzo(a)pyrene	1.067	1.009	5.4	107	-0.02
90	Dibenzo(a,h)anthracene	1.030	0.938	8.9	103	-0.03
91	Benzo(g,h,i)perylene	1.035	0.938	9.4	103	-0.03

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