Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110819\
Data File : BG043567.D
Acq On : 8 Nov 2019 13:32
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 11 00:31:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 04 15:22:15 2019
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	79	0.00
2		1,4-Dioxane	0.425	0.443	-4.2	78	0.00
3		Pyridine	1.073	1.219	-13.6	77	0.00
4		n-Nitrosodimethylamine	0.541	0.582	-7.6	82	0.00
5	S	2-Fluorophenol	1.091	1.176	-7.8	81	0.00
6		Aniline	1.809	1.845	-2.0	75	0.00
7	S	Phenol-d6	1.570	1.659	-5.7	80	0.00
8		2-Chlorophenol	1.205	1.235	-2.5	78	0.00
9		Benzaldehyde	0.772	0.851	-10.2	86	0.00
10	C	Phenol	1.435	1.523	-6.1	79	0.00
11		bis(2-Chloroethyl)ether	1.109	1.036	6.6	70	0.00
12		1,3-Dichlorobenzene	1.510	1.528	-1.2	77	0.00
13	C	1,4-Dichlorobenzene	1.527	1.567	-2.6	78	0.00
14		1,2-Dichlorobenzene	1.491	1.500	-0.6	76	0.00
15		Benzyl Alcohol	1.103	1.127	-2.2	75	0.00
16		2,2'-oxybis(1-Chloropropane	1.613	1.562	3.2	72	-0.01
17		2-Methylphenol	1.046	1.069	-2.2	79	0.00
18		Hexachloroethane	0.551	0.553	-0.4	76	0.00
19	P	n-Nitroso-di-n-propylamine	1.015	0.972	4.2	73	0.00
20		3+4-Methylphenols	1.434	1.406	2.0	75	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	76	0.00
22		Acetophenone	0.512	0.521	-1.8	73	0.00
23	S	Nitrobenzene-d5	0.388	0.398	-2.6	75	0.00
24		Nitrobenzene	0.370	0.381	-3.0	75	0.00
25		Isophorone	0.709	0.687	3.1	70	0.00
26	C	2-Nitrophenol	0.178	0.191	-7.3	79	0.00
27		2,4-Dimethylphenol	0.256	0.276	-7.8	79	0.00
28		bis(2-Chloroethoxy)methane	0.391	0.380	2.8	69	0.00
29	C	2,4-Dichlorophenol	0.328	0.348	-6.1	76	0.00
30		1,2,4-Trichlorobenzene	0.413	0.415	-0.5	74	0.00
31		Naphthalene	1.015	1.046	-3.1	76	0.00
32		Benzoic acid	0.179	0.200	-11.7	90	0.02
33		4-Chloroaniline	0.416	0.424	-1.9	72	0.00
34	C	Hexachlorobutadiene	0.305	0.302	1.0	73	0.00
35		Caprolactam	0.113	0.127	-12.4	79	0.00
36	C	4-Chloro-3-methylphenol	0.357	0.369	-3.4	74	0.00
37		2-Methylnaphthalene	0.779	0.780	-0.1	73	0.00
38		1-Methylnaphthalene	0.741	0.750	-1.2	74	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	76	0.00
40		1,2,4,5-Tetrachlorobenzene	0.740	0.734	0.8	72	0.00
41	P	Hexachlorocyclopentadiene	0.389	0.538	-38.3#	100	0.00
42	S	2,4,6-Tribromophenol	0.381	0.390	-2.4	74	0.00
43	C	2,4,6-Trichlorophenol	0.436	0.443	-1.6	73	0.00
44		2,4,5-Trichlorophenol	0.441	0.434	1.6	71	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110819\
Data File : BG043567.D
Acq On : 8 Nov 2019 13:32
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 11 00:31:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 04 15:22:15 2019
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	S	2-Fluorobiphenyl	1.447	1.450	-0.2	72	0.00
46		1,1'-Biphenyl	1.521	1.540	-1.2	73	0.00
47		2-Chloronaphthalene	1.158	1.179	-1.8	75	0.00
48		2-Nitroaniline	0.346	0.381	-10.1	78	0.00
49		Acenaphthylene	1.802	1.874	-4.0	75	0.00
50		Dimethylphthalate	1.549	1.531	1.2	72	0.00
51		2,6-Dinitrotoluene	0.337	0.354	-5.0	75	0.00
52	C	Acenaphthene	1.099	1.137	-3.5	75	0.00
53		3-Nitroaniline	0.325	0.355	-9.2	78	0.00
54	P	2,4-Dinitrophenol	0.178	0.211	-18.5	86	0.00
55		Dibenzofuran	1.782	1.849	-3.8	75	0.00
56	P	4-Nitrophenol	0.218	0.221	-1.4	72	0.02
57		2,4-Dinitrotoluene	0.481	0.522	-8.5	79	0.00
58		Fluorene	1.494	1.589	-6.4	76	0.00
59		2,3,4,6-Tetrachlorophenol	0.468	0.481	-2.8	71	0.00
60		Diethylphthalate	1.557	1.557	0.0	72	0.00
61		4-Chlorophenyl-phenylether	0.872	0.883	-1.3	74	0.00
62		4-Nitroaniline	0.336	0.385	-14.6	81	0.00
63		Azobenzene	1.260	1.326	-5.2	76	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	84	0.00
65		4,6-Dinitro-2-methylphenol	0.118	0.097	17.8	67	0.00
66	c	n-Nitrosodiphenylamine	0.565	0.541	4.2	76	0.00
67		4-Bromophenyl-phenylether	0.269	0.250	7.1	75	0.00
68		Hexachlorobenzene	0.309	0.293	5.2	77	0.00
69		Atrazine	0.196	0.166	15.3	70	0.00
70	C	Pentachlorophenol	0.141	0.124	12.1	68	0.00
71		Phenanthrene	1.024	1.024	0.0	80	0.00
72		Anthracene	1.025	1.031	-0.6	79	0.00
73		Carbazole	0.928	0.938	-1.1	80	0.00
74		Di-n-butylphthalate	1.126	1.131	-0.4	79	0.00
75	C	Fluoranthene	1.335	1.337	-0.1	78	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	79	0.00
77		Benzidine	0.403	0.375	6.9	66	0.00
78		Pyrene	1.238	1.288	-4.0	79	0.00
79	S	Terphenyl-d14	1.066	1.132	-6.2	79	0.00
80		Butylbenzylphthalate	0.469	0.498	-6.2	80	0.00
81		Benzo(a)anthracene	1.253	1.317	-5.1	80	0.00
82		3,3'-Dichlorobenzidine	0.485	0.491	-1.2	76	0.00
83		Chrysene	1.245	1.288	-3.5	78	0.00
84		Bis(2-ethylhexyl)phthalate	0.666	0.698	-4.8	80	0.00
85	c	Di-n-octyl phthalate	1.130	1.190	-5.3	81	0.00
86		Indeno(1,2,3-cd)pyrene	1.562	1.656	-6.0	81	0.00
87	I	Perylene-d12	1.000	1.000	0.0	81	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110819\
Data File : BG043567.D
Acq On : 8 Nov 2019 13:32
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 11 00:31:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 04 15:22:15 2019
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(b)fluoranthene	1.133	1.140	-0.6	78	0.00
89	Benzo(k)fluoranthene	1.140	1.195	-4.8	81	0.00
90 C	Benzo(a)pyrene	1.082	1.107	-2.3	79	0.00
91	Dibenzo(a,h)anthracene	1.106	1.167	-5.5	81	-0.01
92	Benzo(g,h,i)perylene	1.091	1.130	-3.6	80	0.00

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