Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050819\
Data File : BG040789.D
Acq On : 8 May 2019 14:31
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 08 17:51:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 08 17:46:31 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	118	-0.01
2		1,4-Dioxane	40.000	36.274	9.3	111	0.00
3		Pyridine	40.000	40.071	-0.2	117	-0.02
4		n-Nitrosodimethylamine	40.000	38.924	2.7	120	0.00
5	S	2-Fluorophenol	80.000	79.011	1.2	120	-0.01
6		Aniline	40.000	38.549	3.6	116	-0.02
7	S	Phenol-d6	80.000	81.241	-1.6	123	-0.01
8		2-Chlorophenol	40.000	39.047	2.4	117	-0.02
9		Benzaldehyde	40.000	42.240	-5.6	125	-0.01
10	C	Phenol	40.000	40.574	-1.4	122	-0.01
11		bis(2-Chloroethyl)ether	40.000	38.606	3.5	116	-0.02
12		1,3-Dichlorobenzene	40.000	39.623	0.9	120	-0.02
13	C	1,4-Dichlorobenzene	40.000	39.463	1.3	119	-0.01
14		1,2-Dichlorobenzene	40.000	39.331	1.7	119	-0.02
15		Benzyl Alcohol	40.000	38.153	4.6	114	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	33.333	16.7	101	-0.02
17		2-Methylphenol	40.000	39.340	1.6	120	-0.01
18		Hexachloroethane	40.000	37.072	7.3	113	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	36.262	9.3	112	-0.02
20		3+4-Methylphenols	40.000	39.924	0.2	120	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	116	-0.01
22		Acetophenone	40.000	40.782	-2.0	118	-0.02
23	S	Nitrobenzene-d5	80.000	88.146	-10.2	129	-0.01
24		Nitrobenzene	40.000	42.412	-6.0	126	-0.01
25		Isophorone	40.000	38.702	3.2	114	-0.03
26	C	2-Nitrophenol	40.000	49.192	-23.0#	145	-0.02
27		2,4-Dimethylphenol	40.000	40.082	-0.2	118	-0.02
28		bis(2-Chloroethoxy)methane	40.000	39.196	2.0	114	-0.01
29	C	2,4-Dichlorophenol	40.000	43.937	-9.8	126	-0.02
30		1,2,4-Trichlorobenzene	40.000	43.082	-7.7	125	-0.01
31		Naphthalene	40.000	40.958	-2.4	118	-0.02
32		Benzoic acid	40.000	43.752	-9.4	132	-0.03
33		4-Chloroaniline	40.000	41.129	-2.8	122	-0.02
34	C	Hexachlorobutadiene	40.000	44.062	-10.2	129	-0.02
35		Caprolactam	40.000	40.343	-0.9	117	-0.01
36	C	4-Chloro-3-methylphenol	40.000	42.356	-5.9	127	-0.01
37		2-Methylnaphthalene	40.000	41.859	-4.6	121	-0.01
38		1-Methylnaphthalene	40.000	41.371	-3.4	121	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	124	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	42.174	-5.4	131	-0.02
41	P	Hexachlorocyclopentadiene	40.000	33.133	17.2	104	-0.02
42	S	2,4,6-Tribromophenol	80.000	90.645	-13.3	143	-0.01
43	C	2,4,6-Trichlorophenol	40.000	43.246	-8.1	138	-0.02
44		2,4,5-Trichlorophenol	40.000	44.157	-10.4	139	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050819\
Data File : BG040789.D
Acq On : 8 May 2019 14:31
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 08 17:51:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 08 17:46:31 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	Amount	Calc.	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	80.000	80.995	-1.2	125	-0.02
46		1,1'-Biphenyl	40.000	39.557	1.1	123	-0.02
47		2-Chloronaphthalene	40.000	40.135	-0.3	125	-0.02
48		2-Nitroaniline	40.000	45.367	-13.4	145	-0.01
49		Acenaphthylene	40.000	38.965	2.6	120	-0.01
50		Dimethylphthalate	40.000	39.962	0.1	123	-0.02
51		2,6-Dinitrotoluene	40.000	45.153	-12.9	141	-0.02
52	C	Acenaphthene	40.000	39.051	2.4	124	-0.01
53		3-Nitroaniline	40.000	43.910	-9.8	137	-0.01
54	P	2,4-Dinitrophenol	40.000	35.150	12.1	114	0.00
55		Dibenzofuran	40.000	41.071	-2.7	129	-0.01
56	P	4-Nitrophenol	40.000	36.525	8.7	117	0.00
57		2,4-Dinitrotoluene	40.000	48.486	-21.2	154	-0.01
58		Fluorene	40.000	40.232	-0.6	126	-0.01
59		2,3,4,6-Tetrachlorophenol	40.000	44.576	-11.4	138	-0.01
60		Diethylphthalate	40.000	38.726	3.2	121	-0.02
61		4-Chlorophenyl-phenylether	40.000	42.130	-5.3	132	-0.02
62		4-Nitroaniline	40.000	41.915	-4.8	131	-0.01
63		Azobenzene	40.000	36.345	9.1	113	-0.01
64	I	Phenanthrene-d10	20.000	20.000	0.0	124	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	41.905	-4.8	146	-0.01
66	c	n-Nitrosodiphenylamine	40.000	39.418	1.5	124	-0.01
67		4-Bromophenyl-phenylether	40.000	42.545	-6.4	135	-0.02
68		Hexachlorobenzene	40.000	43.020	-7.6	137	-0.01
69		Atrazine	40.000	38.206	4.5	117	-0.02
70	C	Pentachlorophenol	40.000	40.368	-0.9	126	-0.01
71		Phenanthrene	40.000	40.221	-0.6	125	-0.01
72		Anthracene	40.000	40.137	-0.3	125	-0.02
73		Carbazole	40.000	38.573	3.6	120	-0.01
74		Di-n-butylphthalate	40.000	38.079	4.8	119	-0.03
75	C	Fluoranthene	40.000	40.600	-1.5	127	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	128	-0.03
77		Benzidine	40.000	31.663	20.8	96	-0.01
78		Pyrene	40.000	39.760	0.6	125	-0.01
79	S	Terphenyl-d14	80.000	79.146	1.1	124	-0.02
80		Butylbenzylphthalate	40.000	38.240	4.4	122	-0.02
81		Benzo(a)anthracene	40.000	39.317	1.7	125	-0.02
82		3,3'-Dichlorobenzidine	40.000	37.318	6.7	117	-0.03
83		Chrysene	40.000	39.697	0.8	126	-0.02
84		Bis(2-ethylhexyl)phthalate	40.000	36.443	8.9	116	-0.04
85	c	Di-n-octyl phthalate	40.000	36.462	8.8	116	-0.06
86		Indeno(1,2,3-cd)pyrene	40.000	30.684	23.3	99	-0.08
87	I	Perylene-d12	20.000	20.000	0.0	122	-0.04

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050819\
Data File : BG040789.D
Acq On : 8 May 2019 14:31
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 08 17:51:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 08 17:46:31 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	40.000	40.291	-0.7	122	-0.04
89	Benzo(k)fluoranthene	40.000	42.555	-6.4	130	-0.04
90 C	Benzo(a)pyrene	40.000	40.336	-0.8	122	-0.05
91	Dibenzo(a,h)anthracene	40.000	30.464	23.8	92	-0.09
92	Benzo(g,h,i)perylene	40.000	33.049	17.4	100	-0.10

(#) = Out of Range SPCC's out = 0 CCC's out = 1