Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG021116\
Data File : BG020857.D
Acq On : 11 Feb 2016 15:26
Operator : SJ/UM
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG021116

Quant Time: Feb 11 16:14:12 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 11 15:26:53 2016
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	95	0.00
2		1,4-Dioxane	40.000	38.707	3.2	94	0.00
3		Pyridine	40.000	40.379	-0.9	92	0.00
4		n-Nitrosodimethylamine	40.000	41.604	-4.0	96	0.00
5	S	2-Fluorophenol	80.000	80.012	-0.0	94	0.00
6		Aniline	40.000	40.479	-1.2	95	0.00
7	S	Phenol-d6	80.000	81.432	-1.8	95	0.00
8		2-Chlorophenol	40.000	40.226	-0.6	95	0.00
9		Benzaldehyde	40.000	41.025	-2.6	96	0.00
10	C	Phenol	40.000	40.411	-1.0	95	0.00
11		bis(2-Chloroethyl)ether	40.000	40.601	-1.5	94	0.00
12		1,3-Dichlorobenzene	40.000	40.051	-0.1	95	0.00
13	C	1,4-Dichlorobenzene	40.000	40.875	-2.2	96	0.00
14		1,2-Dichlorobenzene	40.000	40.590	-1.5	95	0.00
15		Benzyl Alcohol	40.000	40.366	-0.9	94	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.428	-1.1	95	0.00
17		2-Methylphenol	40.000	40.761	-1.9	96	0.00
18		Hexachloroethane	40.000	40.526	-1.3	95	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.037	-2.6	97	0.00
20		3+4-Methylphenols	40.000	40.705	-1.8	97	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22		Acetophenone	40.000	40.065	-0.2	97	0.00
23	S	Nitrobenzene-d5	80.000	80.047	-0.1	97	0.00
24		Nitrobenzene	40.000	39.870	0.3	98	0.00
25		Isophorone	40.000	40.288	-0.7	97	0.00
26	C	2-Nitrophenol	40.000	41.187	-3.0	96	0.00
27		2,4-Dimethylphenol	40.000	38.797	3.0	92	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.958	0.1	97	0.00
29	C	2,4-Dichlorophenol	40.000	40.624	-1.6	100	0.00
30		1,2,4-Trichlorobenzene	40.000	39.385	1.5	96	0.00
31		Naphthalene	40.000	39.824	0.4	97	0.00
32		Benzoic acid	40.000	40.761	-1.9	101	0.00
33		4-Chloroaniline	40.000	39.653	0.9	97	0.00
34	C	Hexachlorobutadiene	40.000	40.232	-0.6	98	0.00
35		Caprolactam	40.000	40.423	-1.1	99	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.237	-0.6	99	0.00
37		2-Methylnaphthalene	40.000	39.912	0.2	99	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	101	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	39.781	0.5	99	0.00
40	P	Hexachlorocyclopentadiene	40.000	39.830	0.4	97	0.00
41	S	2,4,6-Tribromophenol	80.000	80.852	-1.1	101	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.593	-1.5	99	0.00
43		2,4,5-Trichlorophenol	40.000	40.440	-1.1	100	0.00
44	S	2-Fluorobiphenyl	80.000	79.865	0.2	99	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG021116\
Data File : BG020857.D
Acq On : 11 Feb 2016 15:26
Operator : SJ/UM
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG021116

Quant Time: Feb 11 16:14:12 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 11 15:26:53 2016
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		1,1'-Biphenyl	40.000	39.941	0.1	100	0.00
46		2-Chloronaphthalene	40.000	39.813	0.5	100	0.00
47		2-Nitroaniline	40.000	40.798	-2.0	101	0.00
48		Acenaphthylene	40.000	40.174	-0.4	101	0.00
49		Dimethylphthalate	40.000	40.206	-0.5	100	0.00
50		2,6-Dinitrotoluene	40.000	40.656	-1.6	101	0.00
51	C	Acenaphthene	40.000	39.948	0.1	100	0.00
52		3-Nitroaniline	40.000	40.097	-0.2	100	0.00
53	P	2,4-Dinitrophenol	40.000	38.574	3.6	103	0.00
54		Dibenzofuran	40.000	40.186	-0.5	102	0.00
55	P	4-Nitrophenol	40.000	42.033	-5.1	102	0.00
56		2,4-Dinitrotoluene	40.000	41.615	-4.0	102	0.00
57		Fluorene	40.000	40.051	-0.1	101	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	40.526	-1.3	101	0.00
59		Diethylphthalate	40.000	40.334	-0.8	102	0.00
60		4-Chlorophenyl-phenylether	40.000	39.473	1.3	101	0.00
61		4-Nitroaniline	40.000	40.543	-1.4	100	0.00
62		Azobenzene	40.000	39.761	0.6	100	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
64		4,6-Dinitro-2-methylphenol	40.000	39.605	1.0	104	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.687	-1.7	102	0.00
66		4-Bromophenyl-phenylether	40.000	40.809	-2.0	102	0.00
67		Hexachlorobenzene	40.000	40.040	-0.1	101	0.00
68		Atrazine	40.000	40.063	-0.2	102	0.00
69	C	Pentachlorophenol	40.000	39.709	0.7	100	0.00
70		Phenanthrene	40.000	39.996	0.0	101	0.00
71		Anthracene	40.000	39.678	0.8	100	0.00
72		Carbazole	40.000	39.722	0.7	100	0.00
73		Di-n-butylphthalate	40.000	40.119	-0.3	101	0.00
74	C	Fluoranthene	40.000	39.723	0.7	100	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
76		Benzidine	40.000	39.189	2.0	97	0.00
77		Pyrene	40.000	40.040	-0.1	102	0.00
78	S	Terphenyl-d14	80.000	79.858	0.2	101	0.00
79		Butylbenzylphthalate	40.000	40.530	-1.3	101	0.00
80		Benzo(a)anthracene	40.000	39.754	0.6	102	0.00
81		3,3'-Dichlorobenzidine	40.000	39.328	1.7	99	0.00
82		Chrysene	40.000	40.075	-0.2	102	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.092	-0.2	101	0.00
84	c	Di-n-octyl phthalate	40.000	40.055	-0.1	100	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	40.364	-0.9	102	0.00
86	I	Perylene-d12	20.000	20.000	0.0	102	0.00
87		Benzo(b)fluoranthene	40.000	40.162	-0.4	102	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG021116\
Data File : BG020857.D
Acq On : 11 Feb 2016 15:26
Operator : SJ/UM
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG021116

Quant Time: Feb 11 16:14:12 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 11 15:26:53 2016
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(k)fluoranthene	40.000	39.458	1.4	101	0.00
89 C	Benzo(a)pyrene	40.000	39.836	0.4	101	0.00
90	Dibenzo(a,h)anthracene	40.000	39.770	0.6	102	0.00
91	Benzo(g,h,i)perylene	40.000	39.783	0.5	101	-0.02

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