Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
Data File : BF085683.D
Acq On : 23 Mar 2016 3:12
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 23 04:02:36 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 18 23:31:39 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	87	-0.02
2		1,4-Dioxane	40.000	40.471	-1.2	86	-0.05
3		Pyridine	40.000	43.817	-9.5	94	-0.06
4		n-Nitrosodimethylamine	40.000	40.580	-1.4	89	-0.05
5	S	2-Fluorophenol	80.000	82.090	-2.6	84	-0.02
6		Aniline	40.000	41.209	-3.0	89	-0.02
7	S	Phenol-d6	80.000	79.919	0.1	85	-0.02
8		2-Chlorophenol	40.000	42.104	-5.3	90	-0.02
9		Benzaldehyde	40.000	53.973	-34.9#	107	-0.02
10	C	Phenol	40.000	43.964	-9.9	86	-0.02
11		bis(2-Chloroethyl)ether	40.000	42.525	-6.3	88	-0.02
12		1,3-Dichlorobenzene	40.000	39.543	1.1	87	-0.03
13	C	1,4-Dichlorobenzene	40.000	38.467	3.8	88	-0.02
14		1,2-Dichlorobenzene	40.000	43.106	-7.8	88	-0.02
15		Benzyl Alcohol	40.000	44.830	-12.1	99	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	40.499	-1.2	85	-0.02
17		2-Methylphenol	40.000	42.575	-6.4	93	-0.01
18		Hexachloroethane	40.000	40.079	-0.2	86	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	44.480	-11.2	93	-0.02
20		3+4-Methylphenols	40.000	38.656	3.4	89	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	92	-0.02
22		Acetophenone	40.000	36.992	7.5	91	-0.02
23	S	Nitrobenzene-d5	80.000	85.713	-7.1	93	-0.02
24		Nitrobenzene	40.000	37.120	7.2	90	-0.02
25		Isophorone	40.000	39.607	1.0	92	-0.02
26	C	2-Nitrophenol	40.000	43.097	-7.7	88	-0.02
27		2,4-Dimethylphenol	40.000	39.002	2.5	94	-0.02
28		bis(2-Chloroethoxy)methane	40.000	46.202	-15.5	99	-0.02
29	C	2,4-Dichlorophenol	40.000	41.689	-4.2	93	-0.02
30		1,2,4-Trichlorobenzene	40.000	38.032	4.9	90	-0.02
31		Naphthalene	40.000	36.897	7.8	90	-0.02
32		Benzoic acid	40.000	30.892	22.8	69	-0.02
33		4-Chloroaniline	40.000	44.750	-11.9	96	-0.02
34	C	Hexachlorobutadiene	40.000	39.317	1.7	87	-0.02
35		Caprolactam	40.000	39.932	0.2	90	-0.02
36	C	4-Chloro-3-methylphenol	40.000	37.961	5.1	84	-0.02
37		2-Methylnaphthalene	40.000	46.388	-16.0	101	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	95	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	34.554	13.6	92	-0.02
40	P	Hexachlorocyclopentadiene	40.000	36.188	9.5	84	-0.02
41	S	2,4,6-Tribromophenol	80.000	85.712	-7.1	103	-0.02
42	C	2,4,6-Trichlorophenol	40.000	36.986	7.5	92	-0.02
43		2,4,5-Trichlorophenol	40.000	40.437	-1.1	96	-0.02
44	S	2-Fluorobiphenyl	80.000	86.955	-8.7	95	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
Data File : BF085683.D
Acq On : 23 Mar 2016 3:12
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 23 04:02:36 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 18 23:31:39 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	35.184	12.0	88	-0.02
46		2-Chloronaphthalene	40.000	39.094	2.3	93	-0.02
47		2-Nitroaniline	40.000	40.730	-1.8	91	-0.02
48		Acenaphthylene	40.000	40.257	-0.6	93	-0.02
49		Dimethylphthalate	40.000	37.051	7.4	92	-0.02
50		2,6-Dinitrotoluene	40.000	37.283	6.8	96	-0.02
51	C	Acenaphthene	40.000	40.572	-1.4	96	-0.02
52		3-Nitroaniline	40.000	35.599	11.0	92	-0.02
53	P	2,4-Dinitrophenol	40.000	30.636	23.4	77	-0.01
54		Dibenzofuran	40.000	45.601	-14.0	107	-0.02
55	P	4-Nitrophenol	40.000	31.648	20.9	72	-0.02
56		2,4-Dinitrotoluene	40.000	44.084	-10.2	98	-0.02
57		Fluorene	40.000	42.151	-5.4	97	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	41.465	-3.7	103	-0.02
59		Diethylphthalate	40.000	40.118	-0.3	99	-0.02
60		4-Chlorophenyl-phenylether	40.000	39.668	0.8	104	-0.02
61		4-Nitroaniline	40.000	40.026	-0.1	88	-0.02
62		Azobenzene	40.000	43.984	-10.0	106	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	93	-0.03
64		4,6-Dinitro-2-methylphenol	40.000	36.120	9.7	83	-0.01
65	c	n-Nitrosodiphenylamine	40.000	41.920	-4.8	94	-0.02
66		4-Bromophenyl-phenylether	40.000	44.695	-11.7	100	-0.02
67		Hexachlorobenzene	40.000	44.095	-10.2	95	-0.02
68		Atrazine	40.000	32.033	19.9	69	-0.02
69	C	Pentachlorophenol	40.000	35.907	10.2	79	-0.02
70		Phenanthrene	40.000	42.342	-5.9	94	-0.02
71		Anthracene	40.000	38.423	3.9	99	-0.02
72		Carbazole	40.000	43.844	-9.6	98	-0.02
73		Di-n-butylphthalate	40.000	41.885	-4.7	94	-0.02
74	C	Fluoranthene	40.000	36.803	8.0	89	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	92	-0.02
76		Benzidine	40.000	26.014	35.0#	59	-0.02
77		Pyrene	40.000	38.304	4.2	85	-0.03
78	S	Terphenyl-d14	80.000	86.209	-7.8	108	-0.02
79		Butylbenzylphthalate	40.000	36.520	8.7	82	-0.03
80		Benzo(a)anthracene	40.000	37.769	5.6	89	-0.02
81		3,3'-Dichlorobenzidine	40.000	39.120	2.2	92	-0.02
82		Chrysene	40.000	35.568	11.1	75	-0.03
83		Bis(2-ethylhexyl)phthalate	40.000	37.169	7.1	86	-0.02
84	c	Di-n-octyl phthalate	40.000	35.365	11.6	79	-0.03
85		Indeno(1,2,3-cd)pyrene	40.000	32.801	18.0	81	-0.06
86	I	Perylene-d12	20.000	20.000	0.0	74	-0.05
87		Benzo(b)fluoranthene	40.000	36.986	7.5	76	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032216\
Data File : BF085683.D
Acq On : 23 Mar 2016 3:12
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 23 04:02:36 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 18 23:31:39 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	48.171	-20.4	74	-0.03
89 C	Benzo(a)pyrene	40.000	41.832	-4.6	76	-0.03
90	Dibenzo(a,h)anthracene	40.000	42.040	-5.1	81	-0.06
91	Benzo(g,h,i)perylene	40.000	43.316	-8.3	84	-0.06

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