Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100115\
Data File : BF081951.D
Acq On : 1 Oct 2015 21:12
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 02 03:19:25 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 28 18:10:42 2015
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	121	-0.03
2		1,4-Dioxane	40.000	39.202	2.0	121	-0.03
3		Pyridine	40.000	39.763	0.6	114	-0.06
4		n-Nitrosodimethylamine	40.000	32.574	18.6	100	-0.05
5	S	2-Fluorophenol	80.000	72.575	9.3	113	-0.03
6		Aniline	40.000	35.363	11.6	111	-0.03
7	S	Phenol-d6	80.000	70.364	12.0	111	-0.03
8		2-Chlorophenol	40.000	36.043	9.9	115	-0.05
9		Benzaldehyde	40.000	31.037	22.4	98	-0.03
10	C	Phenol	40.000	35.990	10.0	109	-0.03
11		bis(2-Chloroethyl)ether	40.000	39.025	2.4	129	-0.03
12		1,3-Dichlorobenzene	40.000	36.332	9.2	115	-0.03
13	C	1,4-Dichlorobenzene	40.000	38.252	4.4	121	-0.03
14		1,2-Dichlorobenzene	40.000	36.682	8.3	121	-0.03
15		Benzyl Alcohol	40.000	32.817	18.0	103	-0.03
16		2,2'-oxybis(1-Chloropropane	40.000	35.049	12.4	110	-0.03
17		2-Methylphenol	40.000	37.605	6.0	117	-0.03
18		Hexachloroethane	40.000	37.612	6.0	120	-0.03
19	P	n-Nitroso-di-n-propylamine	40.000	34.376	14.1	104	-0.03
20		3+4-Methylphenols	40.000	33.661	15.8	108	-0.03
21	I	Naphthalene-d8	20.000	20.000	0.0	121	-0.03
22		Acetophenone	40.000	38.205	4.5	117	-0.03
23	S	Nitrobenzene-d5	80.000	78.427	2.0	122	-0.02
24		Nitrobenzene	40.000	37.846	5.4	113	-0.03
25		Isophorone	40.000	38.294	4.3	118	-0.03
26	C	2-Nitrophenol	40.000	42.066	-5.2	122	-0.03
27		2,4-Dimethylphenol	40.000	38.946	2.6	116	-0.03
28		bis(2-Chloroethoxy)methane	40.000	35.770	10.6	115	-0.03
29	C	2,4-Dichlorophenol	40.000	39.236	1.9	121	-0.03
30		1,2,4-Trichlorobenzene	40.000	39.987	0.0	123	-0.03
31		Naphthalene	40.000	36.496	8.8	119	-0.03
32		Benzoic acid	40.000	34.532	13.7	106	-0.03
33		4-Chloroaniline	40.000	37.670	5.8	113	-0.03
34	C	Hexachlorobutadiene	40.000	40.039	-0.1	123	-0.03
35		Caprolactam	40.000	39.455	1.4	120	-0.03
36	C	4-Chloro-3-methylphenol	40.000	38.503	3.7	117	-0.05
37		2-Methylnaphthalene	40.000	35.931	10.2	109	-0.05
38	I	Acenaphthene-d10	20.000	20.000	0.0	124	-0.03
39		1,2,4,5-Tetrachlorobenzene	40.000	39.594	1.0	128	-0.03
40	P	Hexachlorocyclopentadiene	40.000	38.550	3.6	113	-0.03
41	S	2,4,6-Tribromophenol	80.000	76.985	3.8	126	-0.03
42	C	2,4,6-Trichlorophenol	40.000	35.972	10.1	112	-0.03
43		2,4,5-Trichlorophenol	40.000	42.112	-5.3	131	-0.05
44	S	2-Fluorobiphenyl	80.000	69.164	13.5	107	-0.05

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100115\
Data File : BF081951.D
Acq On : 1 Oct 2015 21:12
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 02 03:19:25 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 28 18:10:42 2015
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	36.791	8.0	120	-0.05
46		2-Chloronaphthalene	40.000	36.299	9.3	113	-0.05
47		2-Nitroaniline	40.000	39.073	2.3	128	-0.03
48		Acenaphthylene	40.000	36.840	7.9	121	-0.03
49		Dimethylphthalate	40.000	38.089	4.8	120	-0.03
50		2,6-Dinitrotoluene	40.000	38.984	2.5	126	-0.03
51	C	Acenaphthene	40.000	37.897	5.3	123	-0.03
52		3-Nitroaniline	40.000	40.219	-0.5	121	-0.05
53	P	2,4-Dinitrophenol	40.000	46.604	-16.5	148	-0.03
54		Dibenzofuran	40.000	38.063	4.8	128	-0.03
55	P	4-Nitrophenol	40.000	34.899	12.8	104	-0.05
56		2,4-Dinitrotoluene	40.000	39.232	1.9	122	-0.05
57		Fluorene	40.000	40.347	-0.9	124	-0.03
58		2,3,4,6-Tetrachlorophenol	40.000	42.289	-5.7	134	-0.03
59		Diethylphthalate	40.000	37.731	5.7	118	-0.05
60		4-Chlorophenyl-phenylether	40.000	39.371	1.6	131	-0.03
61		4-Nitroaniline	40.000	39.492	1.3	128	-0.03
62		Azobenzene	40.000	39.184	2.0	126	-0.03
63	I	Phenanthrene-d10	20.000	20.000	0.0	130	-0.03
64		4,6-Dinitro-2-methylphenol	40.000	44.146	-10.4	136	-0.03
65	c	n-Nitrosodiphenylamine	40.000	34.350	14.1	114	-0.05
66		4-Bromophenyl-phenylether	40.000	40.521	-1.3	138	-0.03
67		Hexachlorobenzene	40.000	34.599	13.5	112	-0.05
68		Atrazine	40.000	34.766	13.1	118	-0.05
69	C	Pentachlorophenol	40.000	39.009	2.5	124	-0.05
70		Phenanthrene	40.000	38.776	3.1	130	-0.03
71		Anthracene	40.000	38.145	4.6	129	-0.05
72		Carbazole	40.000	39.386	1.5	129	-0.05
73		Di-n-butylphthalate	40.000	42.031	-5.1	140	-0.05
74	C	Fluoranthene	40.000	37.237	6.9	133	-0.05
75	I	Chrysene-d12	20.000	20.000	0.0	130	-0.06
76		Benzidine	40.000	43.496	-8.7	128	-0.05
77		Pyrene	40.000	39.749	0.6	124	-0.06
78	S	Terphenyl-d14	80.000	89.749	-12.2	141	-0.05
79		Butylbenzylphthalate	40.000	45.244	-13.1	138	-0.06
80		Benzo(a)anthracene	40.000	37.760	5.6	125	-0.06
81		3,3'-Dichlorobenzidine	40.000	44.937	-12.3	137	-0.05
82		Chrysene	40.000	51.876	-29.7#	144	-0.04
83		Bis(2-ethylhexyl)phthalate	40.000	47.112	-17.8	151	-0.05
84	c	Di-n-octyl phthalate	40.000	48.938	-22.3#	161	-0.05
85		Indeno(1,2,3-cd)pyrene	40.000	37.353	6.6	120	-0.10
86	I	Perylene-d12	20.000	20.000	0.0	129	-0.07
87		Benzo(b)fluoranthene	40.000	34.419	14.0	110	-0.07

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100115\
Data File : BF081951.D
Acq On : 1 Oct 2015 21:12
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 02 03:19:25 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 28 18:10:42 2015
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	46.767	-16.9	169	-0.06
89 C	Benzo(a)pyrene	40.000	39.647	0.9	131	-0.07
90	Dibenzo(a,h)anthracene	40.000	36.720	8.2	122	-0.10
91	Benzo(g,h,i)perylene	40.000	34.717	13.2	117	-0.11

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