Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091216\
Data File : BF090056.D
Acq On : 12 Sep 2016 11:41
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 12 15:36:45 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 12 15:34:24 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	119	-0.02
2		1,4-Dioxane	40.000	39.166	2.1	124	-0.03
3		Pyridine	40.000	36.024	9.9	103	-0.05
4		n-Nitrosodimethylamine	40.000	35.360	11.6	102	-0.03
5	S	2-Fluorophenol	80.000	76.883	3.9	110	-0.02
6		Aniline	40.000	35.036	12.4	110	-0.02
7	S	Phenol-d6	80.000	73.551	8.1	108	-0.02
8		2-Chlorophenol	40.000	40.465	-1.2	119	-0.02
9		Benzaldehyde	40.000	32.330	19.2	96	-0.02
10	C	Phenol	40.000	33.105	17.2	102	-0.02
11		bis(2-Chloroethyl)ether	40.000	39.282	1.8	124	-0.01
12		1,3-Dichlorobenzene	40.000	37.616	6.0	116	-0.02
13	C	1,4-Dichlorobenzene	40.000	38.877	2.8	121	-0.02
14		1,2-Dichlorobenzene	40.000	39.647	0.9	115	-0.02
15		Benzyl Alcohol	40.000	36.831	7.9	110	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	37.936	5.2	111	-0.02
17		2-Methylphenol	40.000	37.880	5.3	113	-0.02
18		Hexachloroethane	40.000	38.815	3.0	116	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	36.127	9.7	103	-0.02
20		3+4-Methylphenols	40.000	38.644	3.4	130	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	121	-0.03
22		Acetophenone	40.000	40.307	-0.8	125	-0.02
23	S	Nitrobenzene-d5	80.000	73.668	7.9	112	-0.02
24		Nitrobenzene	40.000	36.870	7.8	119	-0.02
25		Isophorone	40.000	36.370	9.1	113	-0.02
26	C	2-Nitrophenol	40.000	40.773	-1.9	124	-0.02
27		2,4-Dimethylphenol	40.000	37.971	5.1	126	-0.02
28		bis(2-Chloroethoxy)methane	40.000	36.185	9.5	106	-0.02
29	C	2,4-Dichlorophenol	40.000	39.900	0.3	127	-0.02
30		1,2,4-Trichlorobenzene	40.000	39.741	0.6	117	-0.02
31		Naphthalene	40.000	35.038	12.4	115	-0.02
32		Benzoic acid	40.000	44.906	-12.3	137	-0.01
33		4-Chloroaniline	40.000	38.854	2.9	115	-0.02
34	C	Hexachlorobutadiene	40.000	37.593	6.0	117	-0.02
35		Caprolactam	40.000	36.244	9.4	105	-0.02
36	C	4-Chloro-3-methylphenol	40.000	38.959	2.6	118	-0.02
37		2-Methylnaphthalene	40.000	38.422	3.9	114	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	114	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	38.336	4.2	123	-0.02
40	P	Hexachlorocyclopentadiene	40.000	41.301	-3.3	115	-0.03
41	S	2,4,6-Tribromophenol	80.000	85.001	-6.3	120	-0.02
42	C	2,4,6-Trichlorophenol	40.000	39.675	0.8	107	-0.03
43		2,4,5-Trichlorophenol	40.000	45.525	-13.8	140	-0.02
44	S	2-Fluorobiphenyl	80.000	74.234	7.2	110	-0.02

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	39.405	1.5	124	-0.02
46		2-Chloronaphthalene	40.000	37.965	5.1	113	-0.02
47		2-Nitroaniline	40.000	40.125	-0.3	118	-0.02
48		Acenaphthylene	40.000	38.179	4.6	112	-0.02
49		Dimethylphthalate	40.000	39.674	0.8	128	-0.02
50		2,6-Dinitrotoluene	40.000	44.038	-10.1	129	-0.02
51	C	Acenaphthene	40.000	39.412	1.5	118	-0.02
52		3-Nitroaniline	40.000	38.872	2.8	123	-0.02
53	P	2,4-Dinitrophenol	40.000	44.164	-10.4	145	-0.02
54		Dibenzofuran	40.000	37.557	6.1	110	-0.02
55	P	4-Nitrophenol	40.000	40.139	-0.3	106	-0.02
56		2,4-Dinitrotoluene	40.000	38.591	3.5	99	-0.01
57		Fluorene	40.000	42.325	-5.8	116	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	40.749	-1.9	121	-0.02
59		Diethylphthalate	40.000	38.596	3.5	117	-0.01
60		4-Chlorophenyl-phenylether	40.000	41.832	-4.6	119	-0.02
61		4-Nitroaniline	40.000	42.446	-6.1	112	-0.02
62		Azobenzene	40.000	36.941	7.6	114	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	118	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	39.738	0.7	129	-0.02
65	c	n-Nitrosodiphenylamine	40.000	36.223	9.4	114	-0.02
66		4-Bromophenyl-phenylether	40.000	38.976	2.6	120	-0.02
67		Hexachlorobenzene	40.000	35.762	10.6	119	-0.02
68		Atrazine	40.000	37.125	7.2	119	-0.02
69	C	Pentachlorophenol	40.000	43.081	-7.7	114	-0.02
70		Phenanthrene	40.000	37.782	5.5	114	-0.02
71		Anthracene	40.000	37.350	6.6	125	-0.02
72		Carbazole	40.000	38.663	3.3	120	-0.02
73		Di-n-butylphthalate	40.000	37.309	6.7	118	-0.02
74	C	Fluoranthene	40.000	37.433	6.4	109	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	107	-0.03
76		Benzidine	40.000	38.897	2.8	103	-0.03
77		Pyrene	40.000	37.135	7.2	109	-0.03
78	S	Terphenyl-d14	80.000	76.902	3.9	122	-0.02
79		Butylbenzylphthalate	40.000	43.323	-8.3	119	-0.02
80		Benzo(a)anthracene	40.000	39.551	1.1	108	-0.03
81		3,3'-Dichlorobenzidine	40.000	43.968	-9.9	121	-0.03
82		Chrysene	40.000	42.770	-6.9	121	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	37.750	5.6	101	-0.03
84	c	Di-n-octyl phthalate	40.000	43.097	-7.7	107	-0.03
85		Indeno(1,2,3-cd)pyrene	40.000	41.077	-2.7	110	-0.05
86	I	Perylene-d12	20.000	20.000	0.0	115	-0.03
87		Benzo(b)fluoranthene	40.000	35.248	11.9	89	-0.03

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	40.273	-0.7	136	-0.03
89 C	Benzo(a)pyrene	40.000	39.217	2.0	111	-0.02
90	Dibenzo(a,h)anthracene	40.000	38.398	4.0	110	-0.05
91	Benzo(g,h,i)perylene	40.000	38.173	4.6	111	-0.05

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