Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF112215\
Data File : BF083151.D
Acq On : 22 Nov 2015 11:23
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 23 03:57:43 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Nov 22 09:22:46 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	110	-0.01
2		1,4-Dioxane	40.000	39.551	1.1	114	-0.06
3		Pyridine	40.000	42.598	-6.5	119	-0.27
4		n-Nitrosodimethylamine	40.000	44.820	-12.1	113	-0.08
5	S	2-Fluorophenol	80.000	79.409	0.7	112	-0.03
6		Aniline	40.000	43.597	-9.0	116	-0.02
7	S	Phenol-d6	80.000	79.333	0.8	115	-0.02
8		2-Chlorophenol	40.000	40.566	-1.4	115	-0.01
9		Benzaldehyde	40.000	43.262	-8.2	115	-0.02
10	C	Phenol	40.000	39.780	0.5	123	-0.02
11		bis(2-Chloroethyl)ether	40.000	43.513	-8.8	123	-0.01
12		1,3-Dichlorobenzene	40.000	40.023	-0.1	114	0.00
13	C	1,4-Dichlorobenzene	40.000	39.212	2.0	112	0.00
14		1,2-Dichlorobenzene	40.000	36.959	7.6	102	-0.01
15		Benzyl Alcohol	40.000	37.125	7.2	101	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	44.606	-11.5	127	-0.01
17		2-Methylphenol	40.000	39.245	1.9	112	-0.02
18		Hexachloroethane	40.000	40.156	-0.4	111	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	38.947	2.6	114	-0.02
20		3+4-Methylphenols	40.000	41.157	-2.9	114	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	115	-0.01
22		Acetophenone	40.000	37.947	5.1	115	-0.02
23	S	Nitrobenzene-d5	80.000	77.096	3.6	113	-0.01
24		Nitrobenzene	40.000	39.480	1.3	116	-0.01
25		Isophorone	40.000	39.494	1.3	116	-0.03
26	C	2-Nitrophenol	40.000	39.710	0.7	107	-0.01
27		2,4-Dimethylphenol	40.000	40.710	-1.8	123	-0.01
28		bis(2-Chloroethoxy)methane	40.000	37.887	5.3	108	-0.02
29	C	2,4-Dichlorophenol	40.000	39.397	1.5	113	-0.01
30		1,2,4-Trichlorobenzene	40.000	37.752	5.6	111	-0.01
31		Naphthalene	40.000	39.380	1.5	116	-0.01
32		Benzoic acid	40.000	43.139	-7.8	115	-0.05
33		4-Chloroaniline	40.000	40.857	-2.1	120	-0.01
34	C	Hexachlorobutadiene	40.000	38.120	4.7	112	0.00
35		Caprolactam	40.000	39.267	1.8	116	-0.07
36	C	4-Chloro-3-methylphenol	40.000	39.144	2.1	116	-0.02
37		2-Methylnaphthalene	40.000	36.337	9.2	109	-0.01
38	I	Acenaphthene-d10	20.000	20.000	0.0	116	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	38.243	4.4	115	-0.01
40	P	Hexachlorocyclopentadiene	40.000	32.065	19.8	91	-0.01
41	S	2,4,6-Tribromophenol	80.000	73.839	7.7	105	-0.01
42	C	2,4,6-Trichlorophenol	40.000	37.448	6.4	110	-0.02
43		2,4,5-Trichlorophenol	40.000	38.447	3.9	115	-0.02
44	S	2-Fluorobiphenyl	80.000	73.428	8.2	122	0.00

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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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	36.578	8.6	103	-0.01
46		2-Chloronaphthalene	40.000	37.990	5.0	114	-0.01
47		2-Nitroaniline	40.000	40.224	-0.6	123	-0.01
48		Acenaphthylene	40.000	39.261	1.8	116	-0.01
49		Dimethylphthalate	40.000	38.176	4.6	113	-0.01
50		2,6-Dinitrotoluene	40.000	36.238	9.4	116	-0.01
51	C	Acenaphthene	40.000	37.856	5.4	112	-0.01
52		3-Nitroaniline	40.000	37.413	6.5	106	-0.02
53	P	2,4-Dinitrophenol	40.000	32.848	17.9	94	-0.01
54		Dibenzofuran	40.000	38.041	4.9	111	-0.01
55	P	4-Nitrophenol	40.000	36.170	9.6	108	-0.02
56		2,4-Dinitrotoluene	40.000	35.798	10.5	112	-0.01
57		Fluorene	40.000	36.722	8.2	112	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	38.026	4.9	109	-0.01
59		Diethylphthalate	40.000	35.900	10.3	105	-0.02
60		4-Chlorophenyl-phenylether	40.000	34.963	12.6	108	-0.01
61		4-Nitroaniline	40.000	40.305	-0.8	111	-0.02
62		Azobenzene	40.000	37.616	6.0	114	-0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	110	-0.01
64		4,6-Dinitro-2-methylphenol	40.000	40.844	-2.1	104	-0.01
65	c	n-Nitrosodiphenylamine	40.000	38.079	4.8	111	-0.01
66		4-Bromophenyl-phenylether	40.000	40.114	-0.3	110	-0.01
67		Hexachlorobenzene	40.000	36.394	9.0	103	-0.01
68		Atrazine	40.000	42.304	-5.8	115	-0.02
69	C	Pentachlorophenol	40.000	37.803	5.5	102	-0.01
70		Phenanthrene	40.000	37.103	7.2	104	-0.01
71		Anthracene	40.000	39.502	1.2	118	0.00
72		Carbazole	40.000	40.988	-2.5	112	-0.01
73		Di-n-butylphthalate	40.000	38.821	2.9	110	-0.01
74	C	Fluoranthene	40.000	37.624	5.9	113	-0.01
75	I	Chrysene-d12	20.000	20.000	0.0	106	-0.01
76		Benzidine	40.000	45.048	-12.6	111	-0.01
77		Pyrene	40.000	40.054	-0.1	111	-0.01
78	S	Terphenyl-d14	80.000	80.396	-0.5	108	-0.01
79		Butylbenzylphthalate	40.000	42.240	-5.6	113	-0.01
80		Benzo(a)anthracene	40.000	40.292	-0.7	108	-0.01
81		3,3'-Dichlorobenzidine	40.000	38.046	4.9	103	-0.02
82		Chrysene	40.000	38.517	3.7	109	-0.01
83		Bis(2-ethylhexyl)phthalate	40.000	40.997	-2.5	113	-0.01
84	c	Di-n-octyl phthalate	40.000	40.731	-1.8	112	-0.02
85		Indeno(1,2,3-cd)pyrene	40.000	39.400	1.5	115	-0.02
86	I	Perylene-d12	20.000	20.000	0.0	112	-0.01
87		Benzo(b)fluoranthene	40.000	40.401	-1.0	113	-0.01

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	37.272	6.8	113	-0.01
89 C	Benzo(a)pyrene	40.000	38.110	4.7	110	-0.02
90	Dibenzo(a,h)anthracene	40.000	37.622	5.9	112	-0.02
91	Benzo(g,h,i)perylene	40.000	37.306	6.7	111	-0.02

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