Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070815\
Data File : BF080400.D
Acq On : 8 Jul 2015 21:18
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 09 02:32:24 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 09 00:53:49 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	88	-0.02
2		1,4-Dioxane	40.000	35.386	11.5	77	-0.06
3		Pyridine	40.000	38.002	5.0	88	-0.05
4		n-Nitrosodimethylamine	40.000	40.781	-2.0	90	-0.06
5	S	2-Fluorophenol	80.000	77.240	3.5	86	-0.03
6		Aniline	40.000	43.623	-9.1	96	-0.02
7	S	Phenol-d6	80.000	84.175	-5.2	96	-0.01
8		2-Chlorophenol	40.000	42.388	-6.0	93	-0.02
9		Benzaldehyde	40.000	43.965	-9.9	97	-0.02
10	C	Phenol	40.000	46.256	-15.6	107	-0.02
11		bis(2-Chloroethyl)ether	40.000	39.666	0.8	92	-0.02
12		1,3-Dichlorobenzene	40.000	40.526	-1.3	89	-0.02
13	C	1,4-Dichlorobenzene	40.000	41.098	-2.7	91	-0.02
14		1,2-Dichlorobenzene	40.000	40.692	-1.7	90	-0.02
15		Benzyl Alcohol	40.000	43.532	-8.8	93	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	44.989	-12.5	101	-0.01
17		2-Methylphenol	40.000	46.329	-15.8	103	-0.02
18		Hexachloroethane	40.000	39.667	0.8	86	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	48.857	-22.1	108	-0.02
20		3+4-Methylphenols	40.000	46.518	-16.3	100	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	103	-0.02
22		Acetophenone	40.000	38.862	2.8	96	-0.02
23	S	Nitrobenzene-d5	80.000	83.005	-3.8	101	-0.02
24		Nitrobenzene	40.000	36.946	7.6	96	-0.02
25		Isophorone	40.000	40.999	-2.5	104	-0.02
26	C	2-Nitrophenol	40.000	43.646	-9.1	106	-0.02
27		2,4-Dimethylphenol	40.000	42.232	-5.6	108	-0.02
28		bis(2-Chloroethoxy)methane	40.000	36.823	7.9	93	-0.02
29	C	2,4-Dichlorophenol	40.000	42.826	-7.1	108	-0.02
30		1,2,4-Trichlorobenzene	40.000	39.073	2.3	100	-0.02
31		Naphthalene	40.000	40.489	-1.2	102	-0.02
32		Benzoic acid	40.000	45.138	-12.8	112	-0.01
33		4-Chloroaniline	40.000	39.383	1.5	102	-0.02
34	C	Hexachlorobutadiene	40.000	37.460	6.3	95	-0.02
35		Caprolactam	40.000	45.746	-14.4	116	-0.02
36	C	4-Chloro-3-methylphenol	40.000	42.610	-6.5	110	-0.01
37		2-Methylnaphthalene	40.000	42.102	-5.3	106	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	116	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	37.295	6.8	106	-0.02
40	P	Hexachlorocyclopentadiene	40.000	26.090	34.8#	68	-0.02
41	S	2,4,6-Tribromophenol	80.000	84.754	-5.9	117	-0.02
42	C	2,4,6-Trichlorophenol	40.000	38.791	3.0	106	-0.02
43		2,4,5-Trichlorophenol	40.000	40.123	-0.3	112	-0.02
44	S	2-Fluorobiphenyl	80.000	72.821	9.0	103	-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	38.979	2.6	111	-0.02
46		2-Chloronaphthalene	40.000	39.105	2.2	109	-0.02
47		2-Nitroaniline	40.000	43.897	-9.7	120	-0.02
48		Acenaphthylene	40.000	39.021	2.4	112	-0.02
49		Dimethylphthalate	40.000	39.046	2.4	120	-0.01
50		2,6-Dinitrotoluene	40.000	40.830	-2.1	113	-0.02
51	C	Acenaphthene	40.000	40.242	-0.6	114	-0.02
52		3-Nitroaniline	40.000	41.889	-4.7	113	-0.02
53	P	2,4-Dinitrophenol	40.000	33.426	16.4	92	-0.02
54		Dibenzofuran	40.000	41.281	-3.2	120	-0.02
55	P	4-Nitrophenol	40.000	45.478	-13.7	118	-0.02
56		2,4-Dinitrotoluene	40.000	46.059	-15.1	129	-0.02
57		Fluorene	40.000	40.541	-1.4	113	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	41.952	-4.9	116	-0.02
59		Diethylphthalate	40.000	40.231	-0.6	113	-0.02
60		4-Chlorophenyl-phenylether	40.000	42.132	-5.3	119	-0.02
61		4-Nitroaniline	40.000	45.239	-13.1	118	-0.02
62		Azobenzene	40.000	42.976	-7.4	119	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	122	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	37.743	5.6	115	-0.02
65	c	n-Nitrosodiphenylamine	40.000	40.461	-1.2	121	-0.02
66		4-Bromophenyl-phenylether	40.000	40.003	-0.0	119	-0.02
67		Hexachlorobenzene	40.000	40.139	-0.3	120	-0.02
68		Atrazine	40.000	42.256	-5.6	128	-0.02
69	C	Pentachlorophenol	40.000	33.286	16.8	105	-0.02
70		Phenanthrene	40.000	40.090	-0.2	127	-0.02
71		Anthracene	40.000	39.744	0.6	117	-0.02
72		Carbazole	40.000	41.743	-4.4	128	-0.02
73		Di-n-butylphthalate	40.000	41.058	-2.6	123	-0.03
74	C	Fluoranthene	40.000	40.778	-1.9	125	-0.05
75	I	Chrysene-d12	20.000	20.000	0.0	119	-0.09
76		Benzidine	40.000	46.051	-15.1	130	-0.05
77		Pyrene	40.000	44.425	-11.1	130	-0.05
78	S	Terphenyl-d14	80.000	86.009	-7.5	129	-0.05
79		Butylbenzylphthalate	40.000	49.155	-22.9	134	-0.07
80		Benzo(a)anthracene	40.000	39.827	0.4	116	-0.08
81		3,3'-Dichlorobenzidine	40.000	41.649	-4.1	115	-0.09
82		Chrysene	40.000	39.266	1.8	112	-0.08
83		Bis(2-ethylhexyl)phthalate	40.000	47.725	-19.3	128	-0.09
84	c	Di-n-octyl phthalate	40.000	46.200	-15.5	126	-0.11
85		Indeno(1,2,3-cd)pyrene	40.000	38.177	4.6	108	-0.14
86	I	Perylene-d12	20.000	20.000	0.0	107	-0.11
87		Benzo(b)fluoranthene	40.000	45.200	-13.0	113	-0.11

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	35.645	10.9	100	-0.11
89 C	Benzo(a)pyrene	40.000	40.883	-2.2	107	-0.13
90	Dibenzo(a,h)anthracene	40.000	41.717	-4.3	109	-0.14
91	Benzo(g,h,i)perylene	40.000	40.664	-1.7	106	-0.14

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