Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF062215\
Data File : BF080106.D
Acq On : 22 Jun 2015 22:38
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 23 01:37:52 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 23 00:53:56 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	101	-0.01
2		1,4-Dioxane	0.537	0.501	6.7	94	-0.01
3		Pyridine	1.428	1.318	7.7	93	0.00
4		n-Nitrosodimethylamine	0.679	0.713	-5.0	99	-0.01
5	S	2-Fluorophenol	1.130	1.137	-0.6	99	-0.01
6		Aniline	2.068	2.317	-12.0	108	0.00
7	S	Phenol-d6	1.503	1.582	-5.3	99	-0.01
8		2-Chlorophenol	1.306	1.357	-3.9	101	-0.01
9		Benzaldehyde	0.868	0.904	-4.1	101	0.00
10	C	Phenol	1.641	1.683	-2.6	99	0.00
11		bis(2-Chloroethyl)ether	1.302	1.312	-0.8	98	0.00
12		1,3-Dichlorobenzene	1.569	1.564	0.3	97	0.00
13	C	1,4-Dichlorobenzene	1.492	1.785	-19.6	117	0.00
14		1,2-Dichlorobenzene	1.452	1.547	-6.5	104	0.00
15		Benzyl Alcohol	1.229	1.254	-2.0	98	0.00
16		2,2'-oxybis(1-Chloropropane	1.932	2.295	-18.8	121	0.00
17		2-Methylphenol	1.115	1.191	-6.8	101	0.00
18		Hexachloroethane	0.497	0.554	-11.5	107	0.00
19	P	n-Nitroso-di-n-propylamine	1.044	1.196	-14.6	120	0.00
20		3+4-Methylphenols	1.440	1.627	-13.0	109	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	-0.01
22		Acetophenone	0.466	0.467	-0.2	99	-0.01
23	S	Nitrobenzene-d5	0.344	0.380	-10.5	112	0.00
24		Nitrobenzene	0.355	0.395	-11.3	113	0.00
25		Isophorone	0.691	0.690	0.1	101	0.00
26	C	2-Nitrophenol	0.148	0.184	-24.3#	127	0.00
27		2,4-Dimethylphenol	0.309	0.343	-11.0	118	0.00
28		bis(2-Chloroethoxy)methane	0.406	0.397	2.2	99	0.00
29	C	2,4-Dichlorophenol	0.265	0.315	-18.9	119	0.00
30		1,2,4-Trichlorobenzene	0.301	0.354	-17.6	122	0.00
31		Naphthalene	1.046	1.012	3.3	98	0.00
32		Benzoic acid	0.187	0.171	8.6	93	0.00
33		4-Chloroaniline	0.448	0.410	8.5	98	-0.01
34	C	Hexachlorobutadiene	0.185	0.201	-8.6	117	0.00
35		Caprolactam	0.086	0.091	-5.8	102	-0.01
36	C	4-Chloro-3-methylphenol	0.299	0.293	2.0	97	0.00
37		2-Methylnaphthalene	0.676	0.745	-10.2	111	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	112	0.00
39		1,2,4,5-Tetrachlorobenzene	0.582	0.678	-16.5	125	0.00
40	P	Hexachlorocyclopentadiene	0.260	0.280	-7.7	114	0.00
41	S	2,4,6-Tribromophenol	0.196	0.209	-6.6	115	0.00
42	C	2,4,6-Trichlorophenol	0.396	0.435	-9.8	116	0.00
43		2,4,5-Trichlorophenol	0.456	0.434	4.8	100	-0.01
44	S	2-Fluorobiphenyl	1.402	1.359	3.1	106	0.00

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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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.627	1.664	-2.3	112	0.00
46		2-Chloronaphthalene	1.320	1.251	5.2	101	0.00
47		2-Nitroaniline	0.362	0.364	-0.6	102	0.00
48		Acenaphthylene	2.204	2.179	1.1	102	0.00
49		Dimethylphthalate	1.504	1.567	-4.2	116	0.00
50		2,6-Dinitrotoluene	0.301	0.307	-2.0	102	0.00
51	C	Acenaphthene	1.348	1.272	5.6	101	-0.01
52		3-Nitroaniline	0.348	0.364	-4.6	107	0.00
53	P	2,4-Dinitrophenol	0.106	0.122	-15.1	120	0.00
54		Dibenzofuran	1.913	1.778	7.1	100	-0.01
55	P	4-Nitrophenol	0.278	0.305	-9.7	122	0.00
56		2,4-Dinitrotoluene	0.383	0.428	-11.7	119	0.00
57		Fluorene	1.518	1.599	-5.3	114	0.00
58		2,3,4,6-Tetrachlorophenol	0.385	0.369	4.2	99	-0.01
59		Diethylphthalate	1.514	1.427	5.7	97	0.00
60		4-Chlorophenyl-phenylether	0.729	0.685	6.0	103	-0.01
61		4-Nitroaniline	0.359	0.343	4.5	99	0.00
62		Azobenzene	1.541	1.480	4.0	99	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	118	-0.02
64		4,6-Dinitro-2-methylphenol	0.094	0.094	0.0	108	-0.01
65	c	n-Nitrosodiphenylamine	0.651	0.600	7.8	102	0.00
66		4-Bromophenyl-phenylether	0.213	0.216	-1.4	114	0.00
67		Hexachlorobenzene	0.236	0.238	-0.8	113	-0.01
68		Atrazine	0.206	0.204	1.0	107	-0.01
69	C	Pentachlorophenol	0.134	0.130	3.0	114	-0.01
70		Phenanthrene	1.211	1.156	4.5	110	-0.02
71		Anthracene	1.166	1.124	3.6	112	-0.02
72		Carbazole	1.113	1.008	9.4	102	-0.03
73		Di-n-butylphthalate	1.286	1.077	16.3	100	-0.06
74	C	Fluoranthene	1.290	1.177	8.8	107	-0.11
75	I	Chrysene-d12	1.000	1.000	0.0	110	-0.26
76		Benzidine	0.559	0.608	-8.8	115	-0.11
77		Pyrene	1.231	1.208	1.9	106	-0.13
78	S	Terphenyl-d14	0.856	0.829	3.2	107	-0.15
79		Butylbenzylphthalate	0.495	0.483	2.4	102	-0.21
80		Benzo(a)anthracene	1.218	1.158	4.9	105	-0.26
81		3,3'-Dichlorobenzidine	0.420	0.397	5.5	102	-0.26
82		Chrysene	1.124	1.081	3.8	105	-0.26
83		Bis(2-ethylhexyl)phthalate	0.782	0.740	5.4	100	-0.27
84	c	Di-n-octyl phthalate	1.339	1.214	9.3	97	-0.38
85		Indeno(1,2,3-cd)pyrene	1.417	1.360	4.0	105	-0.41
86	I	Perylene-d12	1.000	1.000	0.0	107	-0.39
87		Benzo(b)fluoranthene	1.211	1.168	3.6	100	-0.38

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.233	1.231	0.2	107	-0.38
89 C	Benzo(a)pyrene	1.150	1.123	2.3	103	-0.39
90	Dibenzo(a,h)anthracene	1.177	1.156	1.8	104	-0.42
91	Benzo(g,h,i)perylene	1.188	1.152	3.0	102	-0.42

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