Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080715\
Data File : BF080939.D
Acq On : 7 Aug 2015 23:04
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 08 06:05:32 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 08 01:33:05 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	102	0.00
2		1,4-Dioxane	0.581	0.562	3.3	97	0.00
3		Pyridine	1.646	1.667	-1.3	107	0.00
4		n-Nitrosodimethylamine	0.841	0.913	-8.6	106	0.01
5	S	2-Fluorophenol	1.224	1.267	-3.5	105	0.01
6		Aniline	2.464	2.438	1.1	102	0.00
7	S	Phenol-d6	1.677	1.768	-5.4	99	0.00
8		2-Chlorophenol	1.416	1.497	-5.7	100	0.00
9		Benzaldehyde	1.131	1.144	-1.1	98	0.00
10	C	Phenol	1.990	2.246	-12.9	100	0.00
11		bis(2-Chloroethyl)ether	1.486	1.428	3.9	96	0.00
12		1,3-Dichlorobenzene	1.503	1.585	-5.5	101	0.00
13	C	1,4-Dichlorobenzene	1.562	1.529	2.1	103	0.00
14		1,2-Dichlorobenzene	1.412	1.436	-1.7	99	0.00
15		Benzyl Alcohol	1.373	1.442	-5.0	98	0.00
16		2,2'-oxybis(1-Chloropropane	2.989	2.866	4.1	100	0.00
17		2-Methylphenol	1.209	1.301	-7.6	106	0.00
18		Hexachloroethane	0.598	0.608	-1.7	102	0.00
19	P	n-Nitroso-di-n-propylamine	1.226	1.246	-1.6	99	0.00
20		3+4-Methylphenols	1.519	1.405	7.5	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
22		Acetophenone	0.490	0.453	7.6	99	0.00
23	S	Nitrobenzene-d5	0.423	0.436	-3.1	101	0.00
24		Nitrobenzene	0.435	0.396	9.0	105	0.00
25		Isophorone	0.813	0.794	2.3	100	0.00
26	C	2-Nitrophenol	0.183	0.184	-0.5	100	0.00
27		2,4-Dimethylphenol	0.343	0.352	-2.6	102	0.00
28		bis(2-Chloroethoxy)methane	0.488	0.423	13.3	99	0.00
29	C	2,4-Dichlorophenol	0.282	0.265	6.0	97	0.00
30		1,2,4-Trichlorobenzene	0.308	0.313	-1.6	99	0.00
31		Naphthalene	1.102	1.110	-0.7	102	0.00
32		Benzoic acid	0.205	0.218	-6.3	118	0.00
33		4-Chloroaniline	0.448	0.456	-1.8	97	0.00
34	C	Hexachlorobutadiene	0.183	0.172	6.0	104	0.00
35		Caprolactam	0.101	0.101	0.0	93	0.00
36	C	4-Chloro-3-methylphenol	0.344	0.327	4.9	96	0.00
37		2-Methylnaphthalene	0.702	0.687	2.1	101	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
39		1,2,4,5-Tetrachlorobenzene	0.617	0.613	0.6	101	0.00
40	P	Hexachlorocyclopentadiene	0.336	0.307	8.6	87	0.00
41	S	2,4,6-Tribromophenol	0.220	0.232	-5.5	96	0.00
42	C	2,4,6-Trichlorophenol	0.452	0.405	10.4	97	0.00
43		2,4,5-Trichlorophenol	0.455	0.476	-4.6	107	0.00
44	S	2-Fluorobiphenyl	1.421	1.387	2.4	101	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.716	1.792	-4.4	101	0.00
46		2-Chloronaphthalene	1.338	1.383	-3.4	101	0.00
47		2-Nitroaniline	0.537	0.579	-7.8	101	0.00
48		Acenaphthylene	2.327	2.429	-4.4	101	0.00
49		Dimethylphthalate	1.661	1.477	11.1	101	0.00
50		2,6-Dinitrotoluene	0.343	0.319	7.0	99	0.00
51	C	Acenaphthene	1.425	1.446	-1.5	98	0.00
52		3-Nitroaniline	0.423	0.406	4.0	94	0.00
53	P	2,4-Dinitrophenol	0.181	0.181	0.0	91	0.00
54		Dibenzofuran	1.927	2.043	-6.0	103	0.00
55	P	4-Nitrophenol	0.317	0.336	-6.0	94	0.00
56		2,4-Dinitrotoluene	0.459	0.447	2.6	105	0.00
57		Fluorene	1.490	1.634	-9.7	105	0.00
58		2,3,4,6-Tetrachlorophenol	0.360	0.357	0.8	99	0.00
59		Diethylphthalate	1.729	1.650	4.6	101	0.00
60		4-Chlorophenyl-phenylether	0.726	0.661	9.0	104	0.00
61		4-Nitroaniline	0.435	0.477	-9.7	105	0.01
62		Azobenzene	1.934	1.873	3.2	100	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
64		4,6-Dinitro-2-methylphenol	0.124	0.128	-3.2	91	0.00
65	c	n-Nitrosodiphenylamine	0.695	0.748	-7.6	101	0.00
66		4-Bromophenyl-phenylether	0.212	0.211	0.5	99	0.00
67		Hexachlorobenzene	0.236	0.234	0.8	100	0.00
68		Atrazine	0.204	0.221	-8.3	102	0.00
69	C	Pentachlorophenol	0.137	0.133	2.9	98	0.00
70		Phenanthrene	1.133	1.139	-0.5	100	0.00
71		Anthracene	1.121	1.111	0.9	99	0.00
72		Carbazole	1.092	1.140	-4.4	99	0.00
73		Di-n-butylphthalate	1.365	1.421	-4.1	106	0.00
74	C	Fluoranthene	1.224	1.250	-2.1	95	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	86	0.00
76		Benzidine	0.788	0.970	-23.1	93	0.00
77		Pyrene	1.477	1.699	-15.0	97	0.00
78	S	Terphenyl-d14	0.968	1.076	-11.2	99	0.00
79		Butylbenzylphthalate	0.712	0.800	-12.4	91	-0.01
80		Benzo(a)anthracene	1.263	1.347	-6.7	90	-0.01
81		3,3'-Dichlorobenzidine	0.460	0.544	-18.3	87	0.00
82		Chrysene	1.210	1.413	-16.8	92	0.00
83		Bis(2-ethylhexyl)phthalate	0.991	1.133	-14.3	91	0.00
84	c	Di-n-octyl phthalate	1.680	2.010	-19.6	92	0.00
85		Indeno(1,2,3-cd)pyrene	1.151	1.194	-3.7	80	0.00
86	I	Perylene-d12	1.000	1.000	0.0	74	0.00
87		Benzo(b)fluoranthene	1.391	1.642	-18.0	81	-0.01

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88	Benzo(k)fluoranthene	1.217	1.046	14.1	72	0.00
89 C	Benzo(a)pyrene	1.195	1.205	-0.8	72	-0.01
90	Dibenzo(a,h)anthracene	1.057	1.186	-12.2	81	0.00
91	Benzo(g,h,i)perylene	1.028	1.141	-11.0	79	0.00

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