Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071015\
Data File : BF080425.D
Acq On : 11 Jul 2015 9:50
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 13 01:05:06 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 11 05:35:30 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	106	-0.01
2		1,4-Dioxane	0.562	0.545	3.0	99	-0.02
3		Pyridine	1.248	1.281	-2.6	109	-0.07
4		n-Nitrosodimethylamine	0.684	0.765	-11.8	115	-0.02
5	S	2-Fluorophenol	1.167	1.240	-6.3	108	-0.01
6		Aniline	1.939	2.122	-9.4	117	0.00
7	S	Phenol-d6	1.555	1.628	-4.7	110	0.00
8		2-Chlorophenol	1.406	1.361	3.2	97	0.00
9		Benzaldehyde	0.860	0.864	-0.5	104	0.00
10	C	Phenol	1.718	1.859	-8.2	115	0.00
11		bis(2-Chloroethyl)ether	1.415	1.396	1.3	104	0.00
12		1,3-Dichlorobenzene	1.590	1.693	-6.5	114	0.00
13	C	1,4-Dichlorobenzene	1.793	1.856	-3.5	109	0.00
14		1,2-Dichlorobenzene	1.699	1.607	5.4	101	0.00
15		Benzyl Alcohol	1.280	1.211	5.4	99	0.00
16		2,2'-oxybis(1-Chloropropane	2.040	2.201	-7.9	112	0.00
17		2-Methylphenol	1.222	1.122	8.2	91	0.00
18		Hexachloroethane	0.553	0.528	4.5	97	0.00
19	P	n-Nitroso-di-n-propylamine	1.064	1.207	-13.4	122	-0.01
20		3+4-Methylphenols	1.536	1.558	-1.4	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
22		Acetophenone	0.479	0.454	5.2	97	0.00
23	S	Nitrobenzene-d5	0.351	0.361	-2.8	111	0.00
24		Nitrobenzene	0.388	0.381	1.8	98	0.00
25		Isophorone	0.628	0.629	-0.2	108	0.00
26	C	2-Nitrophenol	0.186	0.164	11.8	91	0.00
27		2,4-Dimethylphenol	0.293	0.302	-3.1	111	0.00
28		bis(2-Chloroethoxy)methane	0.380	0.392	-3.2	108	-0.01
29	C	2,4-Dichlorophenol	0.271	0.298	-10.0	113	-0.01
30		1,2,4-Trichlorobenzene	0.358	0.327	8.7	97	0.00
31		Naphthalene	1.094	1.084	0.9	110	0.00
32		Benzoic acid	0.206	0.194	5.8	103	-0.01
33		4-Chloroaniline	0.371	0.403	-8.6	110	0.00
34	C	Hexachlorobutadiene	0.208	0.209	-0.5	108	0.00
35		Caprolactam	0.077	0.071	7.8	101	-0.02
36	C	4-Chloro-3-methylphenol	0.287	0.306	-6.6	110	-0.01
37		2-Methylnaphthalene	0.722	0.654	9.4	97	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
39		1,2,4,5-Tetrachlorobenzene	0.669	0.672	-0.4	99	0.00
40	P	Hexachlorocyclopentadiene	0.349	0.214	38.7#	61	0.00
41	S	2,4,6-Tribromophenol	0.195	0.177	9.2	83	0.00
42	C	2,4,6-Trichlorophenol	0.385	0.431	-11.9	105	-0.01
43		2,4,5-Trichlorophenol	0.414	0.475	-14.7	110	0.00
44	S	2-Fluorobiphenyl	1.485	1.515	-2.0	104	0.00

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.743	1.728	0.9	99	0.00
46		2-Chloronaphthalene	1.287	1.271	1.2	96	-0.01
47		2-Nitroaniline	0.380	0.332	12.6	83	0.00
48		Acenaphthylene	2.201	2.314	-5.1	105	0.00
49		Dimethylphthalate	1.512	1.600	-5.8	105	-0.01
50		2,6-Dinitrotoluene	0.339	0.311	8.3	89	0.00
51	C	Acenaphthene	1.399	1.356	3.1	96	0.00
52		3-Nitroaniline	0.331	0.319	3.6	92	0.00
53	P	2,4-Dinitrophenol	0.128	0.060	53.1#	44#	-0.01
54		Dibenzofuran	1.853	1.876	-1.2	102	0.00
55	P	4-Nitrophenol	0.237	0.239	-0.8	98	0.00
56		2,4-Dinitrotoluene	0.397	0.420	-5.8	105	0.00
57		Fluorene	1.561	1.554	0.4	101	0.00
58		2,3,4,6-Tetrachlorophenol	0.363	0.378	-4.1	96	0.00
59		Diethylphthalate	1.480	1.576	-6.5	109	0.00
60		4-Chlorophenyl-phenylether	0.752	0.718	4.5	95	0.00
61		4-Nitroaniline	0.320	0.334	-4.4	102	0.00
62		Azobenzene	1.598	1.529	4.3	94	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
64		4,6-Dinitro-2-methylphenol	0.105	0.060	42.9#	49#	0.00
65	c	n-Nitrosodiphenylamine	0.623	0.647	-3.9	91	-0.01
66		4-Bromophenyl-phenylether	0.213	0.236	-10.8	98	0.00
67		Hexachlorobenzene	0.214	0.215	-0.5	91	-0.01
68		Atrazine	0.194	0.205	-5.7	83	-0.01
69	C	Pentachlorophenol	0.122	0.140	-14.8	95	0.00
70		Phenanthrene	1.202	1.264	-5.2	97	0.00
71		Anthracene	1.113	1.189	-6.8	93	-0.01
72		Carbazole	1.027	1.093	-6.4	93	-0.01
73		Di-n-butylphthalate	1.086	1.268	-16.8	94	0.00
74	C	Fluoranthene	1.246	1.372	-10.1	98	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.02
76		Benzidine	0.391	0.529	-35.3#	134	0.00
77		Pyrene	1.281	1.345	-5.0	108	0.01
78	S	Terphenyl-d14	0.921	0.869	5.6	96	0.00
79		Butylbenzylphthalate	0.486	0.548	-12.8	107	0.01
80		Benzo(a)anthracene	1.260	1.311	-4.0	104	0.02
81		3,3'-Dichlorobenzidine	0.377	0.426	-13.0	109	0.02
82		Chrysene	1.161	1.170	-0.8	100	0.01
83		Bis(2-ethylhexyl)phthalate	0.762	0.844	-10.8	108	0.01
84	c	Di-n-octyl phthalate	1.343	1.491	-11.0	107	0.03
85		Indeno(1,2,3-cd)pyrene	1.442	1.423	1.3	101	0.03
86	I	Perylene-d12	1.000	1.000	0.0	102	0.03
87		Benzo(b)fluoranthene	1.389	1.401	-0.9	103	0.03

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88	Benzo(k)fluoranthene	1.136	1.091	4.0	99	0.03
89 C	Benzo(a)pyrene	1.217	1.177	3.3	101	0.03
90	Dibenzo(a,h)anthracene	1.206	1.133	6.1	99	0.03
91	Benzo(g,h,i)perylene	1.198	1.144	4.5	100	0.05

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