Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012615\
Data File : BF077047.D
Acq On : 26 Jan 2015 10:24
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 26 23:53:12 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 23 15:21:24 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	64	0.05
2		1,4-Dioxane	0.521	0.419	19.6	54	0.02
3		Pyridine	1.348	1.474	-9.3	58	0.02
4		n-Nitrosodimethylamine	0.668	0.633	5.2	60	0.02
5	S	2-Fluorophenol	1.216	1.208	0.7	60	0.05
6		Aniline	2.124	2.100	1.1	59	0.03
7	S	Phenol-d6	1.535	1.500	2.3	59	0.02
8		2-Chlorophenol	1.402	1.373	2.1	59	0.05
9		Benzaldehyde	0.894	0.756	15.4	61	0.05
10	C	Phenol	1.629	1.595	2.1	57	0.02
11		bis(2-Chloroethyl)ether	1.275	1.297	-1.7	63	0.03
12		1,3-Dichlorobenzene	1.595	1.644	-3.1	64	0.03
13	C	1,4-Dichlorobenzene	1.692	1.666	1.5	61	0.03
14		1,2-Dichlorobenzene	1.559	1.569	-0.6	61	0.05
15		Benzyl Alcohol	1.242	1.214	2.3	58	0.03
16		2,2'-oxybis(1-Chloropropane	1.714	1.678	2.1	60	0.05
17		2-Methylphenol	1.146	1.112	3.0	58	0.02
18		Hexachloroethane	0.588	0.592	-0.7	61	0.03
19	P	n-Nitroso-di-n-propylamine	1.074	1.061	1.2	59	0.03
20		3+4-Methylphenols	1.517	1.411	7.0	56	0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	61	0.05
22		Acetophenone	0.490	0.509	-3.9	59	0.03
23	S	Nitrobenzene-d5	0.361	0.377	-4.4	59	0.03
24		Nitrobenzene	0.368	0.391	-6.3	60	0.03
25		Isophorone	0.657	0.687	-4.6	59	0.03
26	C	2-Nitrophenol	0.172	0.174	-1.2	54	0.05
27		2,4-Dimethylphenol	0.284	0.296	-4.2	58	0.02
28		bis(2-Chloroethoxy)methane	0.374	0.401	-7.2	59	0.05
29	C	2,4-Dichlorophenol	0.265	0.287	-8.3	60	0.03
30		1,2,4-Trichlorobenzene	0.294	0.320	-8.8	60	0.03
31		Naphthalene	1.030	1.078	-4.7	59	0.05
32		Benzoic acid	0.158	0.142	10.1	51	0.01
33		4-Chloroaniline	0.416	0.428	-2.9	59	0.05
34	C	Hexachlorobutadiene	0.175	0.187	-6.9	60	0.03
35		Caprolactam	0.078	0.080	-2.6	55	0.01
36	C	4-Chloro-3-methylphenol	0.303	0.315	-4.0	59	0.03
37		2-Methylnaphthalene	0.703	0.743	-5.7	61	0.05
38	I	Acenaphthene-d10	1.000	1.000	0.0	60	0.05
39		1,2,4,5-Tetrachlorobenzene	0.494	0.490	0.8	58	0.05
40	P	Hexachlorocyclopentadiene	0.202	0.241	-19.3	58	0.05
41	S	2,4,6-Tribromophenol	0.162	0.167	-3.1	57	0.03
42	C	2,4,6-Trichlorophenol	0.360	0.360	0.0	56	0.05
43		2,4,5-Trichlorophenol	0.367	0.388	-5.7	59	0.03
44	S	2-Fluorobiphenyl	1.314	1.357	-3.3	59	0.03

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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	1.483	1.547	-4.3	59	0.05
46		2-Chloronaphthalene	1.220	1.260	-3.3	60	0.03
47		2-Nitroaniline	0.370	0.373	-0.8	56	0.05
48		Acenaphthylene	2.058	2.094	-1.7	58	0.03
49		Dimethylphthalate	1.418	1.467	-3.5	60	0.05
50		2,6-Dinitrotoluene	0.283	0.297	-4.9	57	0.05
51	C	Acenaphthene	1.168	1.166	0.2	58	0.05
52		3-Nitroaniline	0.339	0.342	-0.9	55	0.05
53	P	2,4-Dinitrophenol	0.095	0.096	-1.1	54	0.05
54		Dibenzofuran	1.701	1.769	-4.0	60	0.05
55	P	4-Nitrophenol	0.222	0.181	18.5	45#	0.03
56		2,4-Dinitrotoluene	0.384	0.412	-7.3	58	0.03
57		Fluorene	1.441	1.447	-0.4	58	0.03
58		2,3,4,6-Tetrachlorophenol	0.268	0.296	-10.4	62	0.05
59		Diethylphthalate	1.433	1.520	-6.1	61	0.05
60		4-Chlorophenyl-phenylether	0.590	0.626	-6.1	62	0.05
61		4-Nitroaniline	0.328	0.342	-4.3	57	0.03
62		Azobenzene	1.494	1.590	-6.4	61	0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	61	0.03
64		4,6-Dinitro-2-methylphenol	0.104	0.120	-15.4	55	0.03
65	c	n-Nitrosodiphenylamine	0.713	0.751	-5.3	59	0.03
66		4-Bromophenyl-phenylether	0.203	0.220	-8.4	60	0.03
67		Hexachlorobenzene	0.214	0.228	-6.5	62	0.03
68		Atrazine	0.216	0.228	-5.6	56	0.03
69	C	Pentachlorophenol	0.082	0.094	-14.6	63	0.03
70		Phenanthrene	1.247	1.288	-3.3	60	0.03
71		Anthracene	1.216	1.298	-6.7	62	0.02
72		Carbazole	1.180	1.256	-6.4	61	0.03
73		Di-n-butylphthalate	1.365	1.529	-12.0	63	0.02
74	C	Fluoranthene	1.278	1.331	-4.1	58	0.03
75	I	Chrysene-d12	1.000	1.000	0.0	65	0.03
76		Benzidine	0.640	0.635	0.8	65	0.02
77		Pyrene	1.371	1.358	0.9	59	0.03
78	S	Terphenyl-d14	0.869	0.894	-2.9	62	0.03
79		Butylbenzylphthalate	0.621	0.655	-5.5	61	0.03
80		Benzo(a)anthracene	1.255	1.307	-4.1	63	0.03
81		3,3'-Dichlorobenzidine	0.399	0.424	-6.3	64	0.03
82		Chrysene	1.144	1.129	1.3	59	0.02
83		Bis(2-ethylhexyl)phthalate	0.859	0.990	-15.3	70	0.02
84	c	Di-n-octyl phthalate	1.491	1.676	-12.4	67	0.00
85		Indeno(1,2,3-cd)pyrene	1.213	1.252	-3.2	61	-0.08
86	I	Perylene-d12	1.000	1.000	0.0	62	-0.02
87		Benzo(b)fluoranthene	1.347	1.495	-11.0	72	0.00

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88	Benzo(k)fluoranthene	1.177	1.122	4.7	51	0.00
89 C	Benzo(a)pyrene	1.188	1.224	-3.0	60	-0.01
90	Dibenzo(a,h)anthracene	1.090	1.178	-8.1	63	-0.08
91	Benzo(g,h,i)perylene	1.084	1.155	-6.5	62	-0.08

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