Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF101215\
Data File : BF082234.D
Acq On : 12 Oct 2015 22:16
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 13 02:54:36 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 13 02:05:16 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	75	0.00
2		1,4-Dioxane	0.498	0.475	4.6	76	-0.06
3		Pyridine	1.158	1.134	2.1	75	-0.05
4		n-Nitrosodimethylamine	0.491	0.496	-1.0	73	-0.06
5	S	2-Fluorophenol	0.991	0.975	1.6	71	-0.01
6		Aniline	1.663	1.653	0.6	72	-0.01
7	S	Phenol-d6	1.290	1.248	3.3	70	-0.01
8		2-Chlorophenol	1.210	1.190	1.7	76	-0.01
9		Benzaldehyde	0.659	0.664	-0.8	70	-0.01
10	C	Phenol	1.487	1.317	11.4	62	-0.01
11		bis(2-Chloroethyl)ether	1.257	1.152	8.4	68	-0.01
12		1,3-Dichlorobenzene	1.409	1.309	7.1	70	-0.01
13	C	1,4-Dichlorobenzene	1.471	1.347	8.4	68	-0.01
14		1,2-Dichlorobenzene	1.397	1.345	3.7	79	0.00
15		Benzyl Alcohol	0.890	0.938	-5.4	76	0.00
16		2,2'-oxybis(1-Chloropropane	1.751	1.624	7.3	74	0.00
17		2-Methylphenol	0.957	0.961	-0.4	76	0.00
18		Hexachloroethane	0.504	0.450	10.7	69	0.00
19	P	n-Nitroso-di-n-propylamine	0.906	0.857	5.4	74	-0.01
20		3+4-Methylphenols	1.140	1.146	-0.5	78	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	75	0.00
22		Acetophenone	0.396	0.382	3.5	74	-0.01
23	S	Nitrobenzene-d5	0.298	0.294	1.3	72	-0.01
24		Nitrobenzene	0.322	0.309	4.0	79	0.00
25		Isophorone	0.601	0.556	7.5	71	-0.01
26	C	2-Nitrophenol	0.161	0.161	0.0	67	-0.01
27		2,4-Dimethylphenol	0.259	0.272	-5.0	81	0.00
28		bis(2-Chloroethoxy)methane	0.350	0.346	1.1	69	-0.01
29	C	2,4-Dichlorophenol	0.259	0.250	3.5	68	0.00
30		1,2,4-Trichlorobenzene	0.299	0.275	8.0	69	-0.01
31		Naphthalene	0.867	0.818	5.7	73	0.00
32		Benzoic acid	0.174	0.165	5.2	73	-0.01
33		4-Chloroaniline	0.360	0.354	1.7	70	-0.01
34	C	Hexachlorobutadiene	0.186	0.165	11.3	67	-0.01
35		Caprolactam	0.075	0.082	-9.3	77	-0.01
36	C	4-Chloro-3-methylphenol	0.254	0.273	-7.5	81	0.00
37		2-Methylnaphthalene	0.601	0.592	1.5	73	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	71	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.595	0.574	3.5	73	0.00
40	P	Hexachlorocyclopentadiene	0.245	0.210	14.3	58	-0.01
41	S	2,4,6-Tribromophenol	0.188	0.195	-3.7	79	0.00
42	C	2,4,6-Trichlorophenol	0.395	0.400	-1.3	78	0.00
43		2,4,5-Trichlorophenol	0.428	0.421	1.6	72	-0.01
44	S	2-Fluorobiphenyl	1.206	1.199	0.6	68	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.525	1.564	-2.6	80	0.00
46		2-Chloronaphthalene	1.201	1.200	0.1	74	-0.01
47		2-Nitroaniline	0.330	0.368	-11.5	83	0.00
48		Acenaphthylene	1.870	1.822	2.6	71	-0.01
49		Dimethylphthalate	1.408	1.376	2.3	79	0.00
50		2,6-Dinitrotoluene	0.297	0.298	-0.3	69	-0.01
51	C	Acenaphthene	1.123	1.109	1.2	71	-0.01
52		3-Nitroaniline	0.336	0.338	-0.6	71	-0.01
53	P	2,4-Dinitrophenol	0.143	0.135	5.6	63	0.00
54		Dibenzofuran	1.541	1.491	3.2	70	-0.01
55	P	4-Nitrophenol	0.192	0.199	-3.6	69	0.00
56		2,4-Dinitrotoluene	0.371	0.401	-8.1	76	0.00
57		Fluorene	1.266	1.207	4.7	70	-0.01
58		2,3,4,6-Tetrachlorophenol	0.350	0.377	-7.7	84	0.00
59		Diethylphthalate	1.313	1.337	-1.8	76	-0.01
60		4-Chlorophenyl-phenylether	0.580	0.611	-5.3	80	0.00
61		4-Nitroaniline	0.326	0.362	-11.0	77	-0.01
62		Azobenzene	1.285	1.316	-2.4	79	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	79	-0.01
64		4,6-Dinitro-2-methylphenol	0.112	0.105	6.3	67	-0.01
65	c	n-Nitrosodiphenylamine	0.599	0.554	7.5	75	-0.01
66		4-Bromophenyl-phenylether	0.200	0.194	3.0	79	0.00
67		Hexachlorobenzene	0.216	0.205	5.1	76	-0.01
68		Atrazine	0.190	0.204	-7.4	94	0.00
69	C	Pentachlorophenol	0.111	0.108	2.7	70	-0.01
70		Phenanthrene	0.980	0.958	2.2	83	0.00
71		Anthracene	1.010	1.008	0.2	77	-0.01
72		Carbazole	0.922	0.922	0.0	82	0.00
73		Di-n-butylphthalate	1.137	1.088	4.3	77	0.00
74	C	Fluoranthene	1.053	1.054	-0.1	78	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	80	0.00
76		Benzidine	0.580	0.549	5.3	74	0.00
77		Pyrene	1.187	1.172	1.3	83	-0.01
78	S	Terphenyl-d14	0.755	0.683	9.5	74	-0.01
79		Butylbenzylphthalate	0.542	0.542	0.0	85	0.00
80		Benzo(a)anthracene	1.041	1.006	3.4	81	0.00
81		3,3'-Dichlorobenzidine	0.395	0.381	3.5	80	0.00
82		Chrysene	1.096	1.101	-0.5	92	0.00
83		Bis(2-ethylhexyl)phthalate	0.773	0.676	12.5	71	-0.01
84	c	Di-n-octyl phthalate	1.327	1.154	13.0	74	0.00
85		Indeno(1,2,3-cd)pyrene	0.872	0.886	-1.6	92	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	82	0.00
87		Benzo(b)fluoranthene	1.217	1.101	9.5	67	0.00

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88	Benzo(k)fluoranthene	1.023	1.043	-2.0	99	-0.01
89 C	Benzo(a)pyrene	1.046	1.011	3.3	82	0.00
90	Dibenzo(a,h)anthracene	0.857	0.900	-5.0	90	0.00
91	Benzo(g,h,i)perylene	0.832	0.897	-7.8	96	0.00

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