Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102215\
Data File : BF082468.D
Acq On : 23 Oct 2015 2:17
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 23 04:17:48 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 21 13:20:24 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	169#	0.01
2		1,4-Dioxane	0.498	0.457	8.2	165#	0.01
3		Pyridine	1.158	1.232	-6.4	183#	0.01
4		n-Nitrosodimethylamine	0.491	0.517	-5.3	171#	0.01
5	S	2-Fluorophenol	0.991	0.961	3.0	157#	0.01
6		Aniline	1.663	1.509	9.3	148	0.00
7	S	Phenol-d6	1.290	1.200	7.0	152#	0.01
8		2-Chlorophenol	1.210	1.164	3.8	166#	0.01
9		Benzaldehyde	0.659	0.714	-8.3	169#	0.01
10	C	Phenol	1.487	1.466	1.4	155#	0.01
11		bis(2-Chloroethyl)ether	1.257	1.077	14.3	142	0.01
12		1,3-Dichlorobenzene	1.409	1.379	2.1	167#	0.01
13	C	1,4-Dichlorobenzene	1.471	1.401	4.8	160#	0.01
14		1,2-Dichlorobenzene	1.397	1.144	18.1	151#	0.01
15		Benzyl Alcohol	0.890	0.821	7.8	150#	0.00
16		2,2'-oxybis(1-Chloropropane	1.751	1.341	23.4	138	0.00
17		2-Methylphenol	0.957	0.923	3.6	164#	0.01
18		Hexachloroethane	0.504	0.271	46.2#	93	0.00
19	P	n-Nitroso-di-n-propylamine	0.906	0.745	17.8	144	0.00
20		3+4-Methylphenols	1.140	1.157	-1.5	176#	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	149	0.01
22		Acetophenone	0.396	0.430	-8.6	165#	0.00
23	S	Nitrobenzene-d5	0.298	0.301	-1.0	147	0.01
24		Nitrobenzene	0.322	0.343	-6.5	175#	0.01
25		Isophorone	0.601	0.610	-1.5	156#	0.01
26	C	2-Nitrophenol	0.161	0.148	8.1	123	0.01
27		2,4-Dimethylphenol	0.259	0.251	3.1	149	0.00
28		bis(2-Chloroethoxy)methane	0.350	0.363	-3.7	145	0.01
29	C	2,4-Dichlorophenol	0.259	0.274	-5.8	147	0.01
30		1,2,4-Trichlorobenzene	0.299	0.295	1.3	148	0.01
31		Naphthalene	0.867	0.783	9.7	140	0.01
32		Benzoic acid	0.174	0.130	25.3#	114	-0.01
33		4-Chloroaniline	0.360	0.336	6.7	132	0.00
34	C	Hexachlorobutadiene	0.186	0.180	3.2	146	0.01
35		Caprolactam	0.075	0.079	-5.3	146	0.00
36	C	4-Chloro-3-methylphenol	0.254	0.272	-7.1	161#	0.01
37		2-Methylnaphthalene	0.601	0.592	1.5	145	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	134	0.01
39		1,2,4,5-Tetrachlorobenzene	0.595	0.562	5.5	135	0.00
40	P	Hexachlorocyclopentadiene	0.245	0.006#	97.6#	3#	0.00
41	S	2,4,6-Tribromophenol	0.188	0.133	29.3#	101	0.01
42	C	2,4,6-Trichlorophenol	0.395	0.404	-2.3	147	0.01
43		2,4,5-Trichlorophenol	0.428	0.405	5.4	130	0.01
44	S	2-Fluorobiphenyl	1.206	1.218	-1.0	129	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.525	1.553	-1.8	149	0.01
46		2-Chloronaphthalene	1.201	1.211	-0.8	140	0.01
47		2-Nitroaniline	0.330	0.357	-8.2	151#	0.00
48		Acenaphthylene	1.870	1.747	6.6	128	0.01
49		Dimethylphthalate	1.408	1.466	-4.1	158#	0.00
50		2,6-Dinitrotoluene	0.297	0.240	19.2	105	0.00
51	C	Acenaphthene	1.123	0.979	12.8	117	0.01
52		3-Nitroaniline	0.336	0.380	-13.1	150#	0.01
53	P	2,4-Dinitrophenol	0.143	0.014#	90.2#	12#	0.00
54		Dibenzofuran	1.541	1.401	9.1	123	0.01
55	P	4-Nitrophenol	0.192	0.166	13.5	109	0.00
56		2,4-Dinitrotoluene	0.371	0.298	19.7	106	0.00
57		Fluorene	1.266	1.116	11.8	122	0.00
58		2,3,4,6-Tetrachlorophenol	0.350	0.299	14.6	125	0.01
59		Diethylphthalate	1.313	1.310	0.2	140	0.00
60		4-Chlorophenyl-phenylether	0.580	0.507	12.6	125	0.00
61		4-Nitroaniline	0.326	0.369	-13.2	146	0.00
62		Azobenzene	1.285	1.069	16.8	120	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	117	0.01
64		4,6-Dinitro-2-methylphenol	0.112	0.016	85.7#	15#	0.00
65	c	n-Nitrosodiphenylamine	0.599	0.589	1.7	118	0.00
66		4-Bromophenyl-phenylether	0.200	0.187	6.5	113	0.00
67		Hexachlorobenzene	0.216	0.205	5.1	113	0.01
68		Atrazine	0.190	0.123	35.3#	84	0.01
69	C	Pentachlorophenol	0.111	0.101	9.0	96	0.00
70		Phenanthrene	0.980	0.895	8.7	115	0.00
71		Anthracene	1.010	1.028	-1.8	116	0.01
72		Carbazole	0.922	0.900	2.4	118	0.01
73		Di-n-butylphthalate	1.137	1.087	4.4	114	0.00
74	C	Fluoranthene	1.053	0.929	11.8	102	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	110	-0.08
76		Benzidine	0.580	0.670	-15.5	124	0.00
77		Pyrene	1.187	1.089	8.3	105	0.00
78	S	Terphenyl-d14	0.755	0.647	14.3	97	-0.01
79		Butylbenzylphthalate	0.542	0.588	-8.5	127	-0.03
80		Benzo(a)anthracene	1.041	1.055	-1.3	116	-0.08
81		3,3'-Dichlorobenzidine	0.395	0.431	-9.1	123	-0.07
82		Chrysene	1.096	1.063	3.0	122	-0.08
83		Bis(2-ethylhexyl)phthalate	0.773	0.742	4.0	107	-0.08
84	c	Di-n-octyl phthalate	1.327	1.316	0.8	116	-0.16
85		Indeno(1,2,3-cd)pyrene	0.872	0.756	13.3	107	-0.22
86	I	Perylene-d12	1.000	1.000	0.0	110	-0.21
87		Benzo(b)fluoranthene	1.217	1.381	-13.5	112	-0.19

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88	Benzo(k)fluoranthene	1.023	0.834	18.5	107	-0.19
89 C	Benzo(a)pyrene	1.046	1.003	4.1	109	-0.21
90	Dibenzo(a,h)anthracene	0.857	0.794	7.4	107	-0.23
91	Benzo(g,h,i)perylene	0.832	0.760	8.7	109	-0.24

(#) = Out of Range SPCC's out = 2 CCC's out = 0