Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091216\
Data File : BF090056.D
Acq On : 12 Sep 2016 11:41
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 12 15:36:45 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 12 15:34:24 2016
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	119	-0.02
2		1,4-Dioxane	0.592	0.580	2.0	124	-0.03
3		Pyridine	1.579	1.422	9.9	103	-0.05
4		n-Nitrosodimethylamine	0.779	0.689	11.6	102	-0.03
5	S	2-Fluorophenol	1.250	1.201	3.9	110	-0.02
6		Aniline	2.072	1.815	12.4	110	-0.02
7	S	Phenol-d6	1.518	1.396	8.0	108	-0.02
8		2-Chlorophenol	1.356	1.371	-1.1	119	-0.02
9		Benzaldehyde	0.980	0.792	19.2	96	-0.02
10	C	Phenol	1.766	1.461	17.3	102	-0.02
11		bis(2-Chloroethyl)ether	1.336	1.312	1.8	124	-0.01
12		1,3-Dichlorobenzene	1.514	1.424	5.9	116	-0.02
13	C	1,4-Dichlorobenzene	1.549	1.506	2.8	121	-0.02
14		1,2-Dichlorobenzene	1.389	1.376	0.9	115	-0.02
15		Benzyl Alcohol	0.951	0.876	7.9	110	-0.02
16		2,2'-oxybis(1-Chloropropane	2.547	2.416	5.1	111	-0.02
17		2-Methylphenol	1.083	1.025	5.4	113	-0.02
18		Hexachloroethane	0.531	0.515	3.0	116	-0.02
19	P	n-Nitroso-di-n-propylamine	0.989	0.894	9.6	103	-0.02
20		3+4-Methylphenols	1.312	1.268	3.4	130	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	121	-0.03
22		Acetophenone	0.449	0.452	-0.7	125	-0.02
23	S	Nitrobenzene-d5	0.345	0.318	7.8	112	-0.02
24		Nitrobenzene	0.367	0.338	7.9	119	-0.02
25		Isophorone	0.701	0.638	9.0	113	-0.02
26	C	2-Nitrophenol	0.175	0.179	-2.3	124	-0.02
27		2,4-Dimethylphenol	0.324	0.308	4.9	126	-0.02
28		bis(2-Chloroethoxy)methane	0.423	0.383	9.5	106	-0.02
29	C	2,4-Dichlorophenol	0.291	0.290	0.3	127	-0.02
30		1,2,4-Trichlorobenzene	0.318	0.316	0.6	117	-0.02
31		Naphthalene	0.996	0.873	12.3	115	-0.02
32		Benzoic acid	0.218	0.245	-12.4	137	-0.01
33		4-Chloroaniline	0.402	0.390	3.0	115	-0.02
34	C	Hexachlorobutadiene	0.186	0.175	5.9	117	-0.02
35		Caprolactam	0.092	0.083	9.8	105	-0.02
36	C	4-Chloro-3-methylphenol	0.308	0.300	2.6	118	-0.02
37		2-Methylnaphthalene	0.658	0.632	4.0	114	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.586	0.562	4.1	123	-0.02
40	P	Hexachlorocyclopentadiene	0.300	0.310	-3.3	115	-0.03
41	S	2,4,6-Tribromophenol	0.197	0.210	-6.6	120	-0.02
42	C	2,4,6-Trichlorophenol	0.406	0.403	0.7	107	-0.03
43		2,4,5-Trichlorophenol	0.390	0.444	-13.8	140	-0.02
44	S	2-Fluorobiphenyl	1.218	1.131	7.1	110	-0.02

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.610	1.586	1.5	124	-0.02
46		2-Chloronaphthalene	1.139	1.081	5.1	113	-0.02
47		2-Nitroaniline	0.372	0.374	-0.5	118	-0.02
48		Acenaphthylene	1.884	1.798	4.6	112	-0.02
49		Dimethylphthalate	1.446	1.434	0.8	128	-0.02
50		2,6-Dinitrotoluene	0.322	0.354	-9.9	129	-0.02
51	C	Acenaphthene	1.193	1.176	1.4	118	-0.02
52		3-Nitroaniline	0.367	0.356	3.0	123	-0.02
53	P	2,4-Dinitrophenol	0.130	0.183	-40.8#	145	-0.02
54		Dibenzofuran	1.599	1.502	6.1	110	-0.02
55	P	4-Nitrophenol	0.277	0.278	-0.4	106	-0.02
56		2,4-Dinitrotoluene	0.408	0.393	3.7	99	-0.01
57		Fluorene	1.169	1.236	-5.7	116	-0.02
58		2,3,4,6-Tetrachlorophenol	0.334	0.341	-2.1	121	-0.02
59		Diethylphthalate	1.360	1.312	3.5	117	-0.01
60		4-Chlorophenyl-phenylether	0.550	0.575	-4.5	119	-0.02
61		4-Nitroaniline	0.360	0.382	-6.1	112	-0.02
62		Azobenzene	1.358	1.254	7.7	114	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	118	-0.02
64		4,6-Dinitro-2-methylphenol	0.128	0.127	0.8	129	-0.02
65	c	n-Nitrosodiphenylamine	0.601	0.545	9.3	114	-0.02
66		4-Bromophenyl-phenylether	0.201	0.196	2.5	120	-0.02
67		Hexachlorobenzene	0.230	0.206	10.4	119	-0.02
68		Atrazine	0.201	0.186	7.5	119	-0.02
69	C	Pentachlorophenol	0.124	0.133	-7.3	114	-0.02
70		Phenanthrene	0.996	0.941	5.5	114	-0.02
71		Anthracene	1.010	0.943	6.6	125	-0.02
72		Carbazole	0.947	0.916	3.3	120	-0.02
73		Di-n-butylphthalate	1.104	1.030	6.7	118	-0.02
74	C	Fluoranthene	1.044	0.977	6.4	109	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	107	-0.03
76		Benzidine	0.758	0.737	2.8	103	-0.03
77		Pyrene	1.410	1.309	7.2	109	-0.03
78	S	Terphenyl-d14	0.810	0.778	4.0	122	-0.02
79		Butylbenzylphthalate	0.569	0.616	-8.3	119	-0.02
80		Benzo(a)anthracene	1.225	1.211	1.1	108	-0.03
81		3,3'-Dichlorobenzidine	0.442	0.486	-10.0	121	-0.03
82		Chrysene	1.144	1.224	-7.0	121	-0.02
83		Bis(2-ethylhexyl)phthalate	0.787	0.742	5.7	101	-0.03
84	c	Di-n-octyl phthalate	1.270	1.368	-7.7	107	-0.03
85		Indeno(1,2,3-cd)pyrene	0.846	0.869	-2.7	110	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	115	-0.03
87		Benzo(b)fluoranthene	1.253	1.104	11.9	89	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.063	1.070	-0.7	136	-0.03
89 C	Benzo(a)pyrene	1.073	1.052	2.0	111	-0.02
90	Dibenzo(a,h)anthracene	0.787	0.755	4.1	110	-0.05
91	Benzo(g,h,i)perylene	0.792	0.756	4.5	111	-0.05

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