Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092239.D
Acq On : 13 Jan 2017 10:19
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 16 09:36:52 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 06 15:43:51 2017
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	68	-0.05
2		1,4-Dioxane	0.590	0.562	4.7	64	-0.07
3		Pyridine	2.058	1.289	37.4#	41#	-0.09
4		n-Nitrosodimethylamine	0.776	0.557	28.2#	47#	-0.09
5	S	2-Fluorophenol	1.198	1.284	-7.2	76	-0.07
6		Aniline	1.899	1.805	4.9	65	-0.07
7	S	Phenol-d6	1.462	1.354	7.4	62	-0.05
8		2-Chlorophenol	1.362	1.349	1.0	66	-0.05
9		Benzaldehyde	0.904	0.798	11.7	64	-0.05
10	C	Phenol	1.666	1.717	-3.1	65	-0.05
11		bis(2-Chloroethyl)ether	1.326	1.355	-2.2	65	-0.05
12		1,3-Dichlorobenzene	1.455	1.424	2.1	63	-0.05
13	C	1,4-Dichlorobenzene	1.480	1.421	4.0	64	-0.05
14		1,2-Dichlorobenzene	1.285	1.167	9.2	63	-0.05
15		Benzyl Alcohol	1.136	0.953	16.1	56	-0.05
16		2,2'-oxybis(1-Chloropropane	1.975	1.643	16.8	56	-0.05
17		2-Methylphenol	1.096	0.851	22.4	53	-0.05
18		Hexachloroethane	0.549	0.449	18.2	53	-0.05
19	P	n-Nitroso-di-n-propylamine	0.973	0.815	16.2	54	-0.05
20		3+4-Methylphenols	1.307	1.172	10.3	58	-0.05
21	I	Naphthalene-d8	1.000	1.000	0.0	57	-0.05
22		Acetophenone	0.493	0.469	4.9	54	-0.05
23	S	Nitrobenzene-d5	0.360	0.335	6.9	55	-0.05
24		Nitrobenzene	0.383	0.361	5.7	49#	-0.05
25		Isophorone	0.664	0.603	9.2	52	-0.05
26	C	2-Nitrophenol	0.189	0.189	0.0	58	-0.05
27		2,4-Dimethylphenol	0.313	0.289	7.7	48#	-0.05
28		bis(2-Chloroethoxy)methane	0.402	0.339	15.7	47#	-0.05
29	C	2,4-Dichlorophenol	0.265	0.258	2.6	54	-0.05
30		1,2,4-Trichlorobenzene	0.291	0.274	5.8	51	-0.05
31		Naphthalene	0.915	0.935	-2.2	53	-0.05
32		Benzoic acid	0.232	0.208	10.3	48#	-0.03
33		4-Chloroaniline	0.413	0.372	9.9	50	-0.05
34	C	Hexachlorobutadiene	0.153	0.154	-0.7	56	-0.07
35		Caprolactam	0.087	0.080	8.0	51	-0.05
36	C	4-Chloro-3-methylphenol	0.305	0.285	6.6	49#	-0.04
37		2-Methylnaphthalene	0.628	0.552	12.1	55	-0.05
38	I	Acenaphthene-d10	1.000	1.000	0.0	59	-0.05
39		1,2,4,5-Tetrachlorobenzene	0.505	0.496	1.8	52	-0.05
40	P	Hexachlorocyclopentadiene	0.199	0.191	4.0	49#	-0.05
41	S	2,4,6-Tribromophenol	0.166	0.162	2.4	59	-0.05
42	C	2,4,6-Trichlorophenol	0.353	0.326	7.6	49#	-0.04
43		2,4,5-Trichlorophenol	0.364	0.329	9.6	51	-0.04
44	S	2-Fluorobiphenyl	1.042	1.046	-0.4	61	-0.05

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

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.369	1.403	-2.5	53	-0.05
46		2-Chloronaphthalene	1.084	1.077	0.6	56	-0.05
47		2-Nitroaniline	0.386	0.333	13.7	45#	-0.05
48		Acenaphthylene	1.668	1.646	1.3	58	-0.05
49		Dimethylphthalate	1.279	1.197	6.4	52	-0.05
50		2,6-Dinitrotoluene	0.280	0.291	-3.9	61	-0.04
51	C	Acenaphthene	1.131	1.101	2.7	58	-0.05
52		3-Nitroaniline	0.346	0.301	13.0	45#	-0.05
53	P	2,4-Dinitrophenol	0.156	0.156	0.0	51	-0.05
54		Dibenzofuran	1.527	1.418	7.1	59	-0.05
55	P	4-Nitrophenol	0.235	0.189	19.6	43#	-0.04
56		2,4-Dinitrotoluene	0.351	0.341	2.8	58	-0.05
57		Fluorene	1.111	1.149	-3.4	58	-0.05
58		2,3,4,6-Tetrachlorophenol	0.288	0.267	7.3	50	-0.05
59		Diethylphthalate	1.242	1.195	3.8	52	-0.05
60		4-Chlorophenyl-phenylether	0.486	0.460	5.3	56	-0.05
61		4-Nitroaniline	0.359	0.354	1.4	52	-0.05
62		Azobenzene	1.368	1.296	5.3	56	-0.05
63	I	Phenanthrene-d10	1.000	1.000	0.0	58	-0.05
64		4,6-Dinitro-2-methylphenol	0.121	0.134	-10.7	60	-0.04
65	c	n-Nitrosodiphenylamine	0.593	0.577	2.7	57	-0.05
66		4-Bromophenyl-phenylether	0.208	0.216	-3.8	61	-0.05
67		Hexachlorobenzene	0.222	0.216	2.7	55	-0.04
68		Atrazine	0.191	0.181	5.2	55	-0.05
69	C	Pentachlorophenol	0.109	0.086	21.1#	41#	-0.05
70		Phenanthrene	1.084	0.990	8.7	51	-0.05
71		Anthracene	1.086	1.116	-2.8	66	-0.05
72		Carbazole	1.005	1.006	-0.1	63	-0.05
73		Di-n-butylphthalate	1.174	1.190	-1.4	63	-0.04
74	C	Fluoranthene	1.082	1.035	4.3	62	-0.07
75	I	Chrysene-d12	1.000	1.000	0.0	54	-0.05
76		Benzidine	0.580	0.596	-2.8	58	-0.05
77		Pyrene	1.467	1.477	-0.7	58	-0.05
78	S	Terphenyl-d14	0.825	0.947	-14.8	68	-0.05
79		Butylbenzylphthalate	0.667	0.717	-7.5	54	-0.05
80		Benzo(a)anthracene	1.129	1.212	-7.4	58	-0.05
81		3,3'-Dichlorobenzidine	0.428	0.435	-1.6	58	-0.05
82		Chrysene	1.132	1.057	6.6	55	-0.05
83		Bis(2-ethylhexyl)phthalate	0.786	0.885	-12.6	61	-0.05
84	c	Di-n-octyl phthalate	1.456	1.432	1.6	53	-0.05
85		Indeno(1,2,3-cd)pyrene	0.981	0.884	9.9	50#	-0.10
86	I	Perylene-d12	1.000	1.000	0.0	51	-0.05
87		Benzo(b)fluoranthene	1.245	1.459	-17.2	56	-0.05

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.062	0.959	9.7	47#	-0.05
89 C	Benzo(a)pyrene	1.105	1.114	-0.8	50	-0.07
90	Dibenzo(a,h)anthracene	0.908	0.929	-2.3	51	-0.10
91	Benzo(g,h,i)perylene	0.945	0.970	-2.6	51	-0.10

(#) = Out of Range SPCC's out = 0 CCC's out = 1