Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF032317\
Data File : BF093892.D
Acq On : 23 Mar 2017 21:48
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 24 12:59:51 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 23 02:02:22 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	85	-0.01
2		1,4-Dioxane	0.569	0.565	0.7	82	-0.03
3		Pyridine	1.550	1.532	1.2	80	-0.01
4		n-Nitrosodimethylamine	0.638	0.625	2.0	79	-0.02
5	S	2-Fluorophenol	1.184	1.191	-0.6	85	0.00
6		Aniline	2.038	2.027	0.5	81	-0.01
7	S	Phenol-d6	1.465	1.479	-1.0	84	0.00
8		2-Chlorophenol	1.302	1.335	-2.5	84	0.00
9		Benzaldehyde	0.989	0.967	2.2	82	0.00
10	C	Phenol	1.667	1.659	0.5	80	0.00
11		bis(2-Chloroethyl)ether	1.303	1.296	0.5	83	-0.01
12		1,3-Dichlorobenzene	1.460	1.472	-0.8	84	-0.01
13	C	1,4-Dichlorobenzene	1.479	1.481	-0.1	83	-0.01
14		1,2-Dichlorobenzene	1.362	1.376	-1.0	85	-0.01
15		Benzyl Alcohol	1.072	1.089	-1.6	83	0.00
16		2,2'-oxybis(1-Chloropropane	2.007	1.918	4.4	78	0.00
17		2-Methylphenol	1.115	1.120	-0.4	83	0.00
18		Hexachloroethane	0.520	0.526	-1.2	82	-0.01
19	P	n-Nitroso-di-n-propylamine	0.941	0.910	3.3	80	-0.01
20		3+4-Methylphenols	1.350	1.377	-2.0	87	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	84	0.00
22		Acetophenone	0.446	0.450	-0.9	87	-0.01
23	S	Nitrobenzene-d5	0.350	0.350	0.0	83	0.00
24		Nitrobenzene	0.346	0.345	0.3	82	-0.01
25		Isophorone	0.616	0.602	2.3	82	0.00
26	C	2-Nitrophenol	0.165	0.181	-9.7	86	0.00
27		2,4-Dimethylphenol	0.284	0.284	0.0	83	0.00
28		bis(2-Chloroethoxy)methane	0.406	0.397	2.2	81	-0.01
29	C	2,4-Dichlorophenol	0.266	0.269	-1.1	83	0.00
30		1,2,4-Trichlorobenzene	0.274	0.275	-0.4	84	-0.01
31		Naphthalene	0.962	0.959	0.3	84	-0.01
32		Benzoic acid	0.189	0.183	3.2	71	0.00
33		4-Chloroaniline	0.390	0.388	0.5	82	0.00
34	C	Hexachlorobutadiene	0.140	0.141	-0.7	84	-0.01
35		Caprolactam	0.085	0.084	1.2	80	0.00
36	C	4-Chloro-3-methylphenol	0.277	0.275	0.7	82	0.00
37		2-Methylnaphthalene	0.641	0.639	0.3	84	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	84	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.547	0.557	-1.8	86	-0.01
40	P	Hexachlorocyclopentadiene	0.256	0.246	3.9	75	-0.01
41	S	2,4,6-Tribromophenol	0.158	0.160	-1.3	83	0.00
42	C	2,4,6-Trichlorophenol	0.387	0.402	-3.9	86	0.00
43		2,4,5-Trichlorophenol	0.419	0.423	-1.0	81	0.00
44	S	2-Fluorobiphenyl	1.311	1.344	-2.5	86	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.613	1.634	-1.3	85	-0.01
46		2-Chloronaphthalene	1.263	1.266	-0.2	84	0.00
47		2-Nitroaniline	0.375	0.376	-0.3	80	0.00
48		Acenaphthylene	2.099	2.095	0.2	83	-0.01
49		Dimethylphthalate	1.422	1.425	-0.2	84	0.00
50		2,6-Dinitrotoluene	0.326	0.337	-3.4	83	0.00
51	C	Acenaphthene	1.225	1.194	2.5	83	-0.01
52		3-Nitroaniline	0.383	0.366	4.4	78	0.00
53	P	2,4-Dinitrophenol	0.155	0.122	21.3	62	0.00
54		Dibenzofuran	1.667	1.651	1.0	82	-0.01
55	P	4-Nitrophenol	0.300	0.247	17.7	65	0.00
56		2,4-Dinitrotoluene	0.409	0.418	-2.2	81	-0.01
57		Fluorene	1.382	1.313	5.0	85	-0.01
58		2,3,4,6-Tetrachlorophenol	0.287	0.284	1.0	80	0.00
59		Diethylphthalate	1.405	1.404	0.1	83	0.00
60		4-Chlorophenyl-phenylether	0.589	0.572	2.9	86	-0.01
61		4-Nitroaniline	0.367	0.326	11.2	71	0.00
62		Azobenzene	1.329	1.282	3.5	79	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	78	-0.01
64		4,6-Dinitro-2-methylphenol	0.122	0.115	5.7	69	0.00
65	c	n-Nitrosodiphenylamine	0.692	0.716	-3.5	81	-0.01
66		4-Bromophenyl-phenylether	0.197	0.211	-7.1	83	-0.01
67		Hexachlorobenzene	0.199	0.217	-9.0	85	-0.01
68		Atrazine	0.207	0.205	1.0	76	0.00
69	C	Pentachlorophenol	0.124	0.113	8.9	68	0.00
70		Phenanthrene	1.113	1.114	-0.1	79	0.00
71		Anthracene	1.146	1.151	-0.4	79	0.00
72		Carbazole	1.175	1.044	11.1	70	0.00
73		Di-n-butylphthalate	1.222	1.261	-3.2	79	0.00
74	C	Fluoranthene	1.139	0.990	13.1	69	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	57	0.00
76		Benzidine	0.792	0.098	87.6#	7#	-0.01
77		Pyrene	1.451	1.674	-15.4	65	-0.01
78	S	Terphenyl-d14	0.892	1.083	-21.4	70	0.00
79		Butylbenzylphthalate	0.618	0.706	-14.2	61	0.00
80		Benzo(a)anthracene	1.166	1.180	-1.2	57	0.00
81		3,3'-Dichlorobenzidine	0.417	0.398	4.6	51	0.00
82		Chrysene	1.082	1.120	-3.5	59	-0.01
83		Bis(2-ethylhexyl)phthalate	0.844	0.921	-9.1	59	0.00
84	c	Di-n-octyl phthalate	1.410	1.436	-1.8	54	-0.01
85		Indeno(1,2,3-cd)pyrene	0.937	1.363	-45.5#	85	0.00
86	I	Perylene-d12	1.000	1.000	0.0	71	0.00
87		Benzo(b)fluoranthene	1.226	1.187	3.2	67	-0.01

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88	Benzo(k)fluoranthene	1.199	1.119	6.7	64	0.00
89 C	Benzo(a)pyrene	1.129	1.151	-1.9	70	0.00
90	Dibenzo(a,h)anthracene	0.956	1.139	-19.1	85	0.00
91	Benzo(g,h,i)perylene	0.964	1.143	-18.6	85	0.00

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