Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040317\
Data File : BF094145.D
Acq On : 4 Apr 2017 3:50
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 04 05:58:11 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 31 14:38:29 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	95	-0.01
2		1,4-Dioxane	0.569	0.559	1.8	92	-0.04
3		Pyridine	1.550	1.435	7.4	85	-0.03
4		n-Nitrosodimethylamine	0.638	0.625	2.0	89	-0.04
5	S	2-Fluorophenol	1.184	1.105	6.7	89	-0.01
6		Aniline	2.038	1.815	10.9	82	-0.01
7	S	Phenol-d6	1.465	1.300	11.3	84	-0.01
8		2-Chlorophenol	1.302	1.213	6.8	86	-0.01
9		Benzaldehyde	0.989	0.911	7.9	87	-0.01
10	C	Phenol	1.667	1.438	13.7	78	-0.01
11		bis(2-Chloroethyl)ether	1.303	1.285	1.4	92	-0.01
12		1,3-Dichlorobenzene	1.460	1.450	0.7	93	-0.01
13	C	1,4-Dichlorobenzene	1.479	1.474	0.3	93	-0.01
14		1,2-Dichlorobenzene	1.362	1.349	1.0	93	-0.01
15		Benzyl Alcohol	1.072	0.945	11.8	81	-0.01
16		2,2'-oxybis(1-Chloropropane	2.007	1.801	10.3	83	0.00
17		2-Methylphenol	1.115	1.014	9.1	85	-0.01
18		Hexachloroethane	0.520	0.366	29.6#	65	-0.01
19	P	n-Nitroso-di-n-propylamine	0.941	0.899	4.5	89	-0.01
20		3+4-Methylphenols	1.350	1.258	6.8	89	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	88	-0.01
22		Acetophenone	0.446	0.485	-8.7	99	-0.02
23	S	Nitrobenzene-d5	0.350	0.340	2.9	85	-0.01
24		Nitrobenzene	0.346	0.334	3.5	84	-0.02
25		Isophorone	0.616	0.596	3.2	85	-0.02
26	C	2-Nitrophenol	0.165	0.047	71.5#	23#	-0.02
27		2,4-Dimethylphenol	0.284	0.265	6.7	81	-0.01
28		bis(2-Chloroethoxy)methane	0.406	0.414	-2.0	90	-0.01
29	C	2,4-Dichlorophenol	0.266	0.218	18.0	71	-0.01
30		1,2,4-Trichlorobenzene	0.274	0.277	-1.1	89	0.00
31		Naphthalene	0.962	0.954	0.8	88	-0.01
32		Benzoic acid	0.189	0.026	86.2#	11#	0.00
33		4-Chloroaniline	0.390	0.378	3.1	85	-0.01
34	C	Hexachlorobutadiene	0.140	0.144	-2.9	90	0.00
35		Caprolactam	0.085	0.067	21.2	68	-0.02
36	C	4-Chloro-3-methylphenol	0.277	0.215	22.4#	67	-0.01
37		2-Methylnaphthalene	0.641	0.620	3.3	85	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	78	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.547	0.633	-15.7	91	-0.01
40	P	Hexachlorocyclopentadiene	0.256	0.019#	92.6#	5#	-0.02
41	S	2,4,6-Tribromophenol	0.158	0.105	33.5#	51	-0.01
42	C	2,4,6-Trichlorophenol	0.387	0.314	18.9	62	0.00
43		2,4,5-Trichlorophenol	0.419	0.315	24.8	56	-0.01
44	S	2-Fluorobiphenyl	1.311	1.474	-12.4	88	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.613	1.751	-8.6	84	-0.01
46		2-Chloronaphthalene	1.263	1.335	-5.7	83	-0.01
47		2-Nitroaniline	0.375	0.358	4.5	71	-0.01
48		Acenaphthylene	2.099	2.096	0.1	77	-0.02
49		Dimethylphthalate	1.422	1.569	-10.3	86	-0.02
50		2,6-Dinitrotoluene	0.326	0.238	27.0#	55	-0.01
51	C	Acenaphthene	1.225	1.202	1.9	78	-0.01
52		3-Nitroaniline	0.383	0.400	-4.4	80	-0.02
53	P	2,4-Dinitrophenol	0.155	0.000#	100.0#	0#	-9.64#
54		Dibenzofuran	1.667	1.655	0.7	77	-0.01
55	P	4-Nitrophenol	0.300	0.094	68.7#	23#	-0.01
56		2,4-Dinitrotoluene	0.409	0.247	39.6#	44#	-0.02
57		Fluorene	1.382	1.295	6.3	78	-0.02
58		2,3,4,6-Tetrachlorophenol	0.287	0.173	39.7#	46#	-0.01
59		Diethylphthalate	1.405	1.472	-4.8	81	-0.02
60		4-Chlorophenyl-phenylether	0.589	0.582	1.2	81	-0.01
61		4-Nitroaniline	0.367	0.376	-2.5	77	-0.02
62		Azobenzene	1.329	1.235	7.1	71	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	66	-0.01
64		4,6-Dinitro-2-methylphenol	0.122	0.001	99.2#	1#	-0.02
65	c	n-Nitrosodiphenylamine	0.692	0.780	-12.7	75	-0.01
66		4-Bromophenyl-phenylether	0.197	0.221	-12.2	73	0.00
67		Hexachlorobenzene	0.199	0.213	-7.0	70	-0.01
68		Atrazine	0.207	0.193	6.8	60	-0.01
69	C	Pentachlorophenol	0.124	0.032	74.2#	16#	0.00
70		Phenanthrene	1.113	1.144	-2.8	69	-0.01
71		Anthracene	1.146	1.181	-3.1	68	-0.01
72		Carbazole	1.175	1.170	0.4	66	-0.01
73		Di-n-butylphthalate	1.222	1.366	-11.8	72	0.00
74	C	Fluoranthene	1.139	1.161	-1.9	68	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	71	-0.01
76		Benzidine	0.792	0.449	43.3#	39#	-0.01
77		Pyrene	1.451	1.394	3.9	68	-0.01
78	S	Terphenyl-d14	0.892	0.881	1.2	71	-0.01
79		Butylbenzylphthalate	0.618	0.646	-4.5	71	0.00
80		Benzo(a)anthracene	1.166	1.179	-1.1	71	-0.01
81		3,3'-Dichlorobenzidine	0.417	0.396	5.0	64	-0.01
82		Chrysene	1.082	1.097	-1.4	72	-0.01
83		Bis(2-ethylhexyl)phthalate	0.844	0.908	-7.6	73	-0.01
84	c	Di-n-octyl phthalate	1.410	1.596	-13.2	76	0.00
85		Indeno(1,2,3-cd)pyrene	0.937	0.800	14.6	63	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	64	0.00
87		Benzo(b)fluoranthene	1.226	1.409	-14.9	72	-0.01

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88	Benzo(k)fluoranthene	1.199	1.223	-2.0	63	-0.01
89 C	Benzo(a)pyrene	1.129	1.173	-3.9	65	-0.01
90	Dibenzo(a,h)anthracene	0.956	0.946	1.0	63	-0.02
91	Benzo(g,h,i)perylene	0.964	0.920	4.6	62	-0.02

(#) = Out of Range SPCC's out = 2 CCC's out = 3