Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040417\
Data File : BF094169.D
Acq On : 4 Apr 2017 15:57
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 05 00:55:37 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	97	-0.02
2		1,4-Dioxane	0.569	0.558	1.9	93	-0.02
3		Pyridine	1.550	1.539	0.7	92	-0.02
4		n-Nitrosodimethylamine	0.638	0.622	2.5	90	-0.02
5	S	2-Fluorophenol	1.184	1.164	1.7	95	-0.02
6		Aniline	2.038	1.834	10.0	84	-0.02
7	S	Phenol-d6	1.465	1.415	3.4	92	-0.02
8		2-Chlorophenol	1.302	1.307	-0.4	94	-0.02
9		Benzaldehyde	0.989	0.901	8.9	87	-0.02
10	C	Phenol	1.667	1.588	4.7	87	-0.02
11		bis(2-Chloroethyl)ether	1.303	1.298	0.4	95	-0.02
12		1,3-Dichlorobenzene	1.460	1.462	-0.1	95	-0.02
13	C	1,4-Dichlorobenzene	1.479	1.499	-1.4	96	-0.02
14		1,2-Dichlorobenzene	1.362	1.378	-1.2	97	-0.02
15		Benzyl Alcohol	1.072	1.028	4.1	89	-0.02
16		2,2'-oxybis(1-Chloropropane	2.007	1.881	6.3	88	-0.02
17		2-Methylphenol	1.115	1.094	1.9	93	-0.02
18		Hexachloroethane	0.520	0.532	-2.3	95	-0.02
19	P	n-Nitroso-di-n-propylamine	0.941	0.918	2.4	93	-0.02
20		3+4-Methylphenols	1.350	1.411	-4.5	101	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	94	-0.02
22		Acetophenone	0.446	0.477	-7.0	103	-0.02
23	S	Nitrobenzene-d5	0.350	0.360	-2.9	95	-0.02
24		Nitrobenzene	0.346	0.350	-1.2	93	-0.02
25		Isophorone	0.616	0.627	-1.8	95	-0.02
26	C	2-Nitrophenol	0.165	0.187	-13.3	99	-0.02
27		2,4-Dimethylphenol	0.284	0.290	-2.1	95	-0.02
28		bis(2-Chloroethoxy)methane	0.406	0.411	-1.2	94	-0.02
29	C	2,4-Dichlorophenol	0.266	0.278	-4.5	96	-0.02
30		1,2,4-Trichlorobenzene	0.274	0.283	-3.3	96	-0.02
31		Naphthalene	0.962	0.980	-1.9	96	-0.02
32		Benzoic acid	0.189	0.161	14.8	70	-0.02
33		4-Chloroaniline	0.390	0.398	-2.1	95	-0.02
34	C	Hexachlorobutadiene	0.140	0.148	-5.7	98	-0.02
35		Caprolactam	0.085	0.088	-3.5	94	-0.03
36	C	4-Chloro-3-methylphenol	0.277	0.291	-5.1	97	-0.02
37		2-Methylnaphthalene	0.641	0.658	-2.7	96	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	97	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.547	0.566	-3.5	100	-0.02
40	P	Hexachlorocyclopentadiene	0.256	0.246	3.9	86	-0.02
41	S	2,4,6-Tribromophenol	0.158	0.172	-8.9	103	-0.02
42	C	2,4,6-Trichlorophenol	0.387	0.395	-2.1	97	-0.02
43		2,4,5-Trichlorophenol	0.419	0.428	-2.1	94	-0.02
44	S	2-Fluorobiphenyl	1.311	1.342	-2.4	99	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.613	1.624	-0.7	97	-0.02
46		2-Chloronaphthalene	1.263	1.263	0.0	97	-0.02
47		2-Nitroaniline	0.375	0.393	-4.8	97	-0.02
48		Acenaphthylene	2.099	2.095	0.2	96	-0.02
49		Dimethylphthalate	1.422	1.473	-3.6	100	-0.02
50		2,6-Dinitrotoluene	0.326	0.348	-6.7	99	-0.02
51	C	Acenaphthene	1.225	1.219	0.5	97	-0.02
52		3-Nitroaniline	0.383	0.406	-6.0	100	-0.02
53	P	2,4-Dinitrophenol	0.155	0.165	-6.5	97	-0.02
54		Dibenzofuran	1.667	1.636	1.9	94	-0.02
55	P	4-Nitrophenol	0.300	0.282	6.0	85	-0.02
56		2,4-Dinitrotoluene	0.409	0.418	-2.2	93	-0.02
57		Fluorene	1.382	1.342	2.9	100	-0.02
58		2,3,4,6-Tetrachlorophenol	0.287	0.293	-2.1	96	-0.02
59		Diethylphthalate	1.405	1.443	-2.7	99	-0.02
60		4-Chlorophenyl-phenylether	0.589	0.573	2.7	99	-0.02
61		4-Nitroaniline	0.367	0.403	-9.8	102	-0.03
62		Azobenzene	1.329	1.331	-0.2	95	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	98	-0.02
64		4,6-Dinitro-2-methylphenol	0.122	0.131	-7.4	99	-0.02
65	c	n-Nitrosodiphenylamine	0.692	0.687	0.7	98	-0.02
66		4-Bromophenyl-phenylether	0.197	0.203	-3.0	101	-0.02
67		Hexachlorobenzene	0.199	0.208	-4.5	103	-0.02
68		Atrazine	0.207	0.216	-4.3	102	-0.02
69	C	Pentachlorophenol	0.124	0.103	16.9	79	-0.02
70		Phenanthrene	1.113	1.109	0.4	100	-0.02
71		Anthracene	1.146	1.155	-0.8	100	-0.02
72		Carbazole	1.175	1.159	1.4	98	-0.02
73		Di-n-butylphthalate	1.222	1.274	-4.3	101	-0.02
74	C	Fluoranthene	1.139	1.162	-2.0	102	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	101	-0.02
76		Benzidine	0.792	0.693	12.5	86	-0.02
77		Pyrene	1.451	1.451	0.0	100	-0.02
78	S	Terphenyl-d14	0.892	0.893	-0.1	102	-0.02
79		Butylbenzylphthalate	0.618	0.669	-8.3	103	-0.02
80		Benzo(a)anthracene	1.166	1.178	-1.0	101	-0.02
81		3,3'-Dichlorobenzidine	0.417	0.438	-5.0	101	-0.02
82		Chrysene	1.082	1.065	1.6	99	-0.02
83		Bis(2-ethylhexyl)phthalate	0.844	0.894	-5.9	102	-0.02
84	c	Di-n-octyl phthalate	1.410	1.546	-9.6	104	-0.02
85		Indeno(1,2,3-cd)pyrene	0.937	0.955	-1.9	106	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	106	-0.02
87		Benzo(b)fluoranthene	1.226	1.261	-2.9	106	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.199	1.191	0.7	102	-0.02
89 C	Benzo(a)pyrene	1.129	1.154	-2.2	105	-0.02
90	Dibenzo(a,h)anthracene	0.956	0.960	-0.4	106	-0.03
91	Benzo(g,h,i)perylene	0.964	0.961	0.3	106	-0.03

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