Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052316\
Data File : BF087321.D
Acq On : 24 May 2016 3:22
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 24 05:55:39 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 24 01:40:43 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	100	0.00
2		1,4-Dioxane	0.601	0.612	-1.8	103	0.00
3		Pyridine	1.313	1.297	1.2	92	0.00
4		n-Nitrosodimethylamine	1.012	1.109	-9.6	106	-0.01
5	S	2-Fluorophenol	1.138	1.229	-8.0	102	0.00
6		Aniline	1.990	2.036	-2.3	98	0.00
7	S	Phenol-d6	1.538	1.540	-0.1	100	0.00
8		2-Chlorophenol	1.419	1.416	0.2	99	-0.01
9		Benzaldehyde	0.849	0.767	9.7	98	0.00
10	C	Phenol	1.679	1.468	12.6	95	-0.01
11		bis(2-Chloroethyl)ether	1.359	1.336	1.7	99	0.00
12		1,3-Dichlorobenzene	1.548	1.530	1.2	100	0.00
13	C	1,4-Dichlorobenzene	1.589	1.571	1.1	100	0.00
14		1,2-Dichlorobenzene	1.470	1.444	1.8	100	0.00
15		Benzyl Alcohol	1.121	1.152	-2.8	96	0.00
16		2,2'-oxybis(1-Chloropropane	3.058	2.960	3.2	99	0.00
17		2-Methylphenol	1.138	1.127	1.0	99	0.00
18		Hexachloroethane	0.593	0.585	1.3	99	-0.01
19	P	n-Nitroso-di-n-propylamine	1.147	1.143	0.3	101	0.00
20		3+4-Methylphenols	1.375	1.429	-3.9	98	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.478	0.482	-0.8	99	0.00
23	S	Nitrobenzene-d5	0.397	0.401	-1.0	97	0.00
24		Nitrobenzene	0.419	0.430	-2.6	98	0.00
25		Isophorone	0.712	0.719	-1.0	99	0.00
26	C	2-Nitrophenol	0.171	0.182	-6.4	98	-0.01
27		2,4-Dimethylphenol	0.297	0.299	-0.7	98	-0.01
28		bis(2-Chloroethoxy)methane	0.401	0.400	0.2	100	0.00
29	C	2,4-Dichlorophenol	0.268	0.280	-4.5	99	0.00
30		1,2,4-Trichlorobenzene	0.307	0.308	-0.3	99	0.00
31		Naphthalene	1.002	0.983	1.9	99	0.00
32		Benzoic acid	0.141	0.138	2.1	89	0.00
33		4-Chloroaniline	0.369	0.368	0.3	96	0.00
34	C	Hexachlorobutadiene	0.186	0.187	-0.5	99	0.00
35		Caprolactam	0.078	0.085	-9.0	103	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.319	-5.3	99	0.00
37		2-Methylnaphthalene	0.626	0.624	0.3	100	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
39		1,2,4,5-Tetrachlorobenzene	0.640	0.638	0.3	98	0.00
40	P	Hexachlorocyclopentadiene	0.206	0.214	-3.9	91	-0.01
41	S	2,4,6-Tribromophenol	0.192	0.195	-1.6	97	0.00
42	C	2,4,6-Trichlorophenol	0.370	0.384	-3.8	98	0.00
43		2,4,5-Trichlorophenol	0.376	0.387	-2.9	100	0.00
44	S	2-Fluorobiphenyl	1.419	1.334	6.0	95	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.684	1.621	3.7	98	0.00
46		2-Chloronaphthalene	1.288	1.273	1.2	99	0.00
47		2-Nitroaniline	0.464	0.482	-3.9	99	0.00
48		Acenaphthylene	2.040	1.991	2.4	98	0.00
49		Dimethylphthalate	1.602	1.591	0.7	98	0.00
50		2,6-Dinitrotoluene	0.323	0.333	-3.1	101	0.00
51	C	Acenaphthene	1.254	1.221	2.6	100	0.00
52		3-Nitroaniline	0.331	0.332	-0.3	97	0.00
53	P	2,4-Dinitrophenol	0.138	0.149	-8.0	91	0.00
54		Dibenzofuran	1.697	1.674	1.4	100	0.00
55	P	4-Nitrophenol	0.158	0.168	-6.3	102	0.00
56		2,4-Dinitrotoluene	0.431	0.443	-2.8	97	-0.01
57		Fluorene	1.472	1.431	2.8	97	-0.01
58		2,3,4,6-Tetrachlorophenol	0.289	0.309	-6.9	99	0.00
59		Diethylphthalate	1.558	1.494	4.1	97	0.00
60		4-Chlorophenyl-phenylether	0.718	0.715	0.4	100	0.00
61		4-Nitroaniline	0.309	0.333	-7.8	96	0.00
62		Azobenzene	1.616	1.663	-2.9	97	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	0.127	0.136	-7.1	97	0.00
65	c	n-Nitrosodiphenylamine	0.647	0.649	-0.3	99	0.00
66		4-Bromophenyl-phenylether	0.219	0.217	0.9	99	-0.01
67		Hexachlorobenzene	0.229	0.226	1.3	98	0.00
68		Atrazine	0.206	0.210	-1.9	101	0.00
69	C	Pentachlorophenol	0.061	0.066	-8.2	96	0.00
70		Phenanthrene	1.108	1.073	3.2	99	0.00
71		Anthracene	1.119	1.090	2.6	99	0.00
72		Carbazole	0.955	0.949	0.6	100	0.00
73		Di-n-butylphthalate	1.234	1.274	-3.2	99	0.00
74	C	Fluoranthene	1.111	1.073	3.4	96	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
76		Benzidine	0.515	0.542	-5.2	94	0.00
77		Pyrene	1.440	1.456	-1.1	99	0.00
78	S	Terphenyl-d14	0.941	0.985	-4.7	101	0.00
79		Butylbenzylphthalate	0.638	0.667	-4.5	95	0.00
80		Benzo(a)anthracene	1.138	1.121	1.5	96	0.00
81		3,3'-Dichlorobenzidine	0.628	0.610	2.9	98	0.00
82		Chrysene	1.129	1.087	3.7	97	0.00
83		Bis(2-ethylhexyl)phthalate	0.932	0.941	-1.0	97	0.00
84	c	Di-n-octyl phthalate	1.421	1.472	-3.6	94	-0.01
85		Indeno(1,2,3-cd)pyrene	0.905	0.931	-2.9	99	0.00
86	I	Perylene-d12	1.000	1.000	0.0	95	0.00
87		Benzo(b)fluoranthene	1.274	1.093	14.2	76	0.00

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88	Benzo(k)fluoranthene	1.097	1.240	-13.0	126	0.00
89 C	Benzo(a)pyrene	1.102	1.123	-1.9	99	0.00
90	Dibenzo(a,h)anthracene	0.940	0.990	-5.3	99	0.00
91	Benzo(g,h,i)perylene	0.927	0.982	-5.9	99	0.00

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