Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050516\
Data File : BF086707.D
Acq On : 5 May 2016 14:04
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 06 01:22:01 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 04 15:12:53 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	1.237	1.171	5.3	100	0.00
3		Pyridine	3.018	3.083	-2.2	100	0.00
4		n-Nitrosodimethylamine	1.855	1.892	-2.0	100	0.00
5	S	2-Fluorophenol	2.415	2.535	-5.0	100	0.00
6		Aniline	3.996	3.918	2.0	100	0.00
7	S	Phenol-d6	3.233	3.136	3.0	100	0.00
8		2-Chlorophenol	2.883	2.840	1.5	100	0.00
9		Benzaldehyde	1.644	1.530	6.9	100	0.00
10	C	Phenol	3.575	3.734	-4.4	100	0.00
11		bis(2-Chloroethyl)ether	3.045	2.766	9.2	100	0.00
12		1,3-Dichlorobenzene	3.073	2.949	4.0	100	0.00
13	C	1,4-Dichlorobenzene	3.172	2.983	6.0	100	0.00
14		1,2-Dichlorobenzene	2.765	2.900	-4.9	100	0.00
15		Benzyl Alcohol	2.093	2.147	-2.6	100	0.00
16		2,2'-oxybis(1-Chloropropane	6.255	5.866	6.2	100	0.00
17		2-Methylphenol	2.336	2.276	2.6	100	0.00
18		Hexachloroethane	1.199	1.197	0.2	100	0.00
19	P	n-Nitroso-di-n-propylamine	2.216	1.976	10.8	100	0.00
20		3+4-Methylphenols	2.700	2.547	5.7	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
22		Acetophenone	0.935	0.980	-4.8	100	0.00
23	S	Nitrobenzene-d5	0.771	0.735	4.7	100	0.00
24		Nitrobenzene	0.810	0.815	-0.6	100	0.00
25		Isophorone	1.445	1.409	2.5	100	0.00
26	C	2-Nitrophenol	0.356	0.347	2.5	100	0.00
27		2,4-Dimethylphenol	0.628	0.596	5.1	100	0.00
28		bis(2-Chloroethoxy)methane	0.802	0.777	3.1	100	0.00
29	C	2,4-Dichlorophenol	0.545	0.557	-2.2	100	0.00
30		1,2,4-Trichlorobenzene	0.605	0.616	-1.8	100	0.00
31		Naphthalene	2.043	1.935	5.3	100	0.00
32		Benzoic acid	0.204	0.227	-11.3	100	0.00
33		4-Chloroaniline	0.785	0.746	5.0	100	0.00
34	C	Hexachlorobutadiene	0.343	0.339	1.2	100	0.00
35		Caprolactam	0.171	0.193	-12.9	100	0.00
36	C	4-Chloro-3-methylphenol	0.610	0.588	3.6	100	0.00
37		2-Methylnaphthalene	1.197	1.180	1.4	100	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
39		1,2,4,5-Tetrachlorobenzene	1.160	1.160	0.0	100	0.00
40	P	Hexachlorocyclopentadiene	0.471	0.482	-2.3	100	0.00
41	S	2,4,6-Tribromophenol	0.391	0.418	-6.9	100	0.00
42	C	2,4,6-Trichlorophenol	0.740	0.759	-2.6	100	0.00
43		2,4,5-Trichlorophenol	0.764	0.775	-1.4	100	0.00
44	S	2-Fluorobiphenyl	2.397	2.187	8.8	100	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	3.264	3.492	-7.0	100	0.00
46		2-Chloronaphthalene	2.523	2.565	-1.7	100	0.00
47		2-Nitroaniline	0.897	0.864	3.7	100	0.00
48		Acenaphthylene	3.735	3.943	-5.6	100	0.00
49		Dimethylphthalate	3.011	3.044	-1.1	100	0.00
50		2,6-Dinitrotoluene	0.639	0.668	-4.5	100	0.00
51	C	Acenaphthene	2.455	2.372	3.4	100	0.00
52		3-Nitroaniline	0.726	0.702	3.3	100	0.00
53	P	2,4-Dinitrophenol	0.172	0.186	-8.1	100	0.00
54		Dibenzofuran	3.058	3.144	-2.8	100	0.00
55	P	4-Nitrophenol	0.423	0.441	-4.3	100	0.00
56		2,4-Dinitrotoluene	0.784	0.781	0.4	100	0.00
57		Fluorene	2.484	2.650	-6.7	100	0.00
58		2,3,4,6-Tetrachlorophenol	0.562	0.560	0.4	100	0.00
59		Diethylphthalate	2.842	2.850	-0.3	100	0.00
60		4-Chlorophenyl-phenylether	1.175	1.114	5.2	100	0.00
61		4-Nitroaniline	0.701	0.735	-4.9	100	0.00
62		Azobenzene	3.118	3.022	3.1	100	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	0.193	0.227	-17.6	100	0.00
65	c	n-Nitrosodiphenylamine	1.269	1.318	-3.9	100	0.00
66		4-Bromophenyl-phenylether	0.397	0.419	-5.5	100	0.00
67		Hexachlorobenzene	0.486	0.484	0.4	100	0.00
68		Atrazine	0.388	0.370	4.6	100	0.00
69	C	Pentachlorophenol	0.207	0.220	-6.3	100	0.00
70		Phenanthrene	2.151	2.080	3.3	100	0.00
71		Anthracene	2.248	2.335	-3.9	100	0.00
72		Carbazole	1.974	2.185	-10.7	100	0.00
73		Di-n-butylphthalate	2.331	2.327	0.2	100	0.00
74	C	Fluoranthene	2.153	2.281	-5.9	100	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
76		Benzidine	0.906	0.871	3.9	100	0.00
77		Pyrene	3.092	2.912	5.8	100	0.00
78	S	Terphenyl-d14	1.904	1.606	15.7	100	0.00
79		Butylbenzylphthalate	1.330	1.387	-4.3	100	0.00
80		Benzo(a)anthracene	2.235	2.124	5.0	100	0.00
81		3,3'-Dichlorobenzidine	1.392	1.596	-14.7	100	0.00
82		Chrysene	2.368	2.255	4.8	100	0.00
83		Bis(2-ethylhexyl)phthalate	1.540	1.634	-6.1	100	0.00
84	c	Di-n-octyl phthalate	2.315	2.707	-16.9	100	0.00
85		Indeno(1,2,3-cd)pyrene	1.988	1.706	14.2	100	0.00
86	I	Perylene-d12	1.000	1.000	0.0	100	0.00
87		Benzo(b)fluoranthene	2.379	2.364	0.6	100	0.00

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88	Benzo(k)fluoranthene	2.349	2.493	-6.1	100	0.00
89 C	Benzo(a)pyrene	2.213	2.196	0.8	100	0.00
90	Dibenzo(a,h)anthracene	2.049	1.726	15.8	100	0.00
91	Benzo(g,h,i)perylene	2.137	1.712	19.9	100	0.00

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