Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122117\
Data File : BG031694.D
Acq On : 22 Dec 2017 4:06
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 22 05:42:12 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 22 03:38:35 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	98	-0.01
2		1,4-Dioxane	0.407	0.364	10.6	91	-0.01
3		Pyridine	1.088	1.085	0.3	93	-0.01
4		n-Nitrosodimethylamine	0.439	0.422	3.9	92	0.00
5	S	2-Fluorophenol	1.098	1.061	3.4	95	0.00
6		Aniline	1.834	1.763	3.9	92	0.00
7	S	Phenol-d6	1.426	1.429	-0.2	95	0.00
8		2-Chlorophenol	1.274	1.259	1.2	96	-0.01
9		Benzaldehyde	0.858	0.824	4.0	92	-0.01
10	C	Phenol	1.439	1.417	1.5	95	0.00
11		bis(2-Chloroethyl)ether	1.195	1.116	6.6	91	0.00
12		1,3-Dichlorobenzene	1.582	1.497	5.4	93	-0.01
13	C	1,4-Dichlorobenzene	1.616	1.531	5.3	94	-0.01
14		1,2-Dichlorobenzene	1.531	1.452	5.2	93	0.00
15		Benzyl Alcohol	1.070	1.067	0.3	94	0.00
16		2,2'-oxybis(1-Chloropropane	0.973	0.870	10.6	88	0.00
17		2-Methylphenol	1.030	1.021	0.9	95	0.00
18		Hexachloroethane	0.489	0.475	2.9	98	-0.01
19	P	n-Nitroso-di-n-propylamine	0.974	0.930	4.5	93	-0.01
20		3+4-Methylphenols	1.464	1.464	0.0	95	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
22		Acetophenone	0.465	0.438	5.8	93	0.00
23	S	Nitrobenzene-d5	0.265	0.269	-1.5	96	-0.01
24		Nitrobenzene	0.273	0.273	0.0	94	0.00
25		Isophorone	0.570	0.537	5.8	91	-0.01
26	C	2-Nitrophenol	0.143	0.155	-8.4	105	-0.01
27		2,4-Dimethylphenol	0.301	0.293	2.7	98	0.00
28		bis(2-Chloroethoxy)methane	0.383	0.354	7.6	91	-0.01
29	C	2,4-Dichlorophenol	0.354	0.353	0.3	96	0.00
30		1,2,4-Trichlorobenzene	0.432	0.408	5.6	93	-0.01
31		Naphthalene	1.009	0.971	3.8	95	-0.01
32		Benzoic acid	0.186	0.195	-4.8	97	0.00
33		4-Chloroaniline	0.449	0.434	3.3	95	0.00
34	C	Hexachlorobutadiene	0.296	0.280	5.4	94	0.00
35		Caprolactam	0.107	0.103	3.7	92	0.00
36	C	4-Chloro-3-methylphenol	0.327	0.334	-2.1	100	0.00
37		2-Methylnaphthalene	0.818	0.780	4.6	94	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
39		1,2,4,5-Tetrachlorobenzene	0.842	0.799	5.1	96	-0.01
40	P	Hexachlorocyclopentadiene	0.444	0.438	1.4	97	0.00
41	S	2,4,6-Tribromophenol	0.305	0.323	-5.9	105	0.00
42	C	2,4,6-Trichlorophenol	0.486	0.489	-0.6	100	0.00
43		2,4,5-Trichlorophenol	0.538	0.538	0.0	100	-0.01
44	S	2-Fluorobiphenyl	1.593	1.475	7.4	95	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.708	1.595	6.6	94	0.00
46		2-Chloronaphthalene	1.303	1.228	5.8	96	-0.01
47		2-Nitroaniline	0.225	0.243	-8.0	107	0.00
48		Acenaphthylene	2.084	1.980	5.0	97	0.00
49		Dimethylphthalate	1.762	1.659	5.8	96	-0.01
50		2,6-Dinitrotoluene	0.324	0.341	-5.2	103	0.00
51	C	Acenaphthene	1.365	1.309	4.1	97	0.00
52		3-Nitroaniline	0.316	0.329	-4.1	102	0.00
53	P	2,4-Dinitrophenol	0.115	0.139	-20.9	124	0.00
54		Dibenzofuran	2.093	1.982	5.3	97	0.00
55	P	4-Nitrophenol	0.257	0.260	-1.2	98	0.00
56		2,4-Dinitrotoluene	0.417	0.463	-11.0	107	0.00
57		Fluorene	1.700	1.594	6.2	97	0.00
58		2,3,4,6-Tetrachlorophenol	0.525	0.525	0.0	101	0.00
59		Diethylphthalate	1.717	1.605	6.5	95	0.00
60		4-Chlorophenyl-phenylether	1.040	0.981	5.7	96	0.00
61		4-Nitroaniline	0.309	0.329	-6.5	100	0.00
62		Azobenzene	1.101	1.013	8.0	93	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	0.086	0.105	-22.1	121	0.00
65	c	n-Nitrosodiphenylamine	0.664	0.633	4.7	96	0.00
66		4-Bromophenyl-phenylether	0.289	0.278	3.8	97	0.00
67		Hexachlorobenzene	0.296	0.294	0.7	100	-0.01
68		Atrazine	0.258	0.246	4.7	95	0.00
69	C	Pentachlorophenol	0.203	0.214	-5.4	101	0.00
70		Phenanthrene	1.132	1.067	5.7	95	0.00
71		Anthracene	1.142	1.078	5.6	96	0.00
72		Carbazole	1.011	0.948	6.2	94	0.00
73		Di-n-butylphthalate	1.123	1.070	4.7	96	-0.01
74	C	Fluoranthene	1.389	1.299	6.5	96	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
76		Benzidine	0.617	0.463	25.0	74	0.00
77		Pyrene	1.278	1.193	6.7	95	0.00
78	S	Terphenyl-d14	0.875	0.824	5.8	97	0.00
79		Butylbenzylphthalate	0.429	0.417	2.8	98	-0.01
80		Benzo(a)anthracene	1.272	1.199	5.7	97	0.00
81		3,3'-Dichlorobenzidine	0.493	0.469	4.9	96	-0.01
82		Chrysene	1.204	1.129	6.2	96	0.00
83		Bis(2-ethylhexyl)phthalate	0.621	0.592	4.7	96	-0.02
84	c	Di-n-octyl phthalate	1.046	1.006	3.8	96	-0.02
85		Indeno(1,2,3-cd)pyrene	1.544	1.528	1.0	99	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	102	-0.01
87		Benzo(b)fluoranthene	1.241	1.172	5.6	99	-0.02

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88	Benzo(k)fluoranthene	1.242	1.167	6.0	96	-0.02
89 C	Benzo(a)pyrene	1.189	1.120	5.8	97	-0.01
90	Dibenzo(a,h)anthracene	1.228	1.182	3.7	99	-0.02
91	Benzo(g,h,i)perylene	1.450	1.402	3.3	99	-0.02

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