Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG122217\
Data File : BG031714.D
Acq On : 22 Dec 2017 16:54
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 23 01:16:52 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 11:08:55 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	82	-0.01
2		1,4-Dioxane	0.407	0.390	4.2	82	0.00
3		Pyridine	1.088	1.108	-1.8	79	-0.01
4		n-Nitrosodimethylamine	0.439	0.430	2.1	79	0.00
5	S	2-Fluorophenol	1.098	1.085	1.2	81	0.00
6		Aniline	1.834	1.729	5.7	75	0.00
7	S	Phenol-d6	1.426	1.421	0.4	79	0.00
8		2-Chlorophenol	1.274	1.259	1.2	80	-0.01
9		Benzaldehyde	0.858	0.824	4.0	77	-0.01
10	C	Phenol	1.439	1.424	1.0	80	0.00
11		bis(2-Chloroethyl)ether	1.195	1.106	7.4	76	0.00
12		1,3-Dichlorobenzene	1.582	1.513	4.4	78	-0.01
13	C	1,4-Dichlorobenzene	1.616	1.525	5.6	78	-0.01
14		1,2-Dichlorobenzene	1.531	1.497	2.2	80	0.00
15		Benzyl Alcohol	1.070	1.046	2.2	77	-0.01
16		2,2'-oxybis(1-Chloropropane	0.973	0.834	14.3	70	0.00
17		2-Methylphenol	1.030	0.998	3.1	78	0.00
18		Hexachloroethane	0.489	0.481	1.6	83	-0.01
19	P	n-Nitroso-di-n-propylamine	0.974	0.900	7.6	75	-0.01
20		3+4-Methylphenols	1.464	1.416	3.3	77	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	81	-0.01
22		Acetophenone	0.465	0.446	4.1	78	-0.01
23	S	Nitrobenzene-d5	0.265	0.270	-1.9	80	-0.01
24		Nitrobenzene	0.273	0.274	-0.4	78	-0.01
25		Isophorone	0.570	0.535	6.1	75	-0.01
26	C	2-Nitrophenol	0.143	0.158	-10.5	88	-0.01
27		2,4-Dimethylphenol	0.301	0.289	4.0	79	-0.01
28		bis(2-Chloroethoxy)methane	0.383	0.358	6.5	76	-0.01
29	C	2,4-Dichlorophenol	0.354	0.348	1.7	78	0.00
30		1,2,4-Trichlorobenzene	0.432	0.410	5.1	77	-0.01
31		Naphthalene	1.009	0.960	4.9	77	-0.01
32		Benzoic acid	0.186	0.190	-2.2	78	-0.02
33		4-Chloroaniline	0.449	0.423	5.8	76	0.00
34	C	Hexachlorobutadiene	0.296	0.294	0.7	81	0.00
35		Caprolactam	0.107	0.104	2.8	76	-0.01
36	C	4-Chloro-3-methylphenol	0.327	0.333	-1.8	82	-0.01
37		2-Methylnaphthalene	0.818	0.780	4.6	78	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	81	0.00
39		1,2,4,5-Tetrachlorobenzene	0.842	0.807	4.2	79	-0.01
40	P	Hexachlorocyclopentadiene	0.444	0.435	2.0	78	-0.01
41	S	2,4,6-Tribromophenol	0.305	0.328	-7.5	86	0.00
42	C	2,4,6-Trichlorophenol	0.486	0.493	-1.4	82	0.00
43		2,4,5-Trichlorophenol	0.538	0.542	-0.7	81	-0.01
44	S	2-Fluorobiphenyl	1.593	1.514	5.0	79	-0.01

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.708	1.627	4.7	78	-0.01
46		2-Chloronaphthalene	1.303	1.245	4.5	79	-0.01
47		2-Nitroaniline	0.225	0.246	-9.3	87	0.00
48		Acenaphthylene	2.084	2.002	3.9	79	0.00
49		Dimethylphthalate	1.762	1.689	4.1	79	-0.01
50		2,6-Dinitrotoluene	0.324	0.358	-10.5	88	0.00
51	C	Acenaphthene	1.365	1.316	3.6	79	0.00
52		3-Nitroaniline	0.316	0.336	-6.3	84	0.00
53	P	2,4-Dinitrophenol	0.115	0.113	1.7	82	-0.01
54		Dibenzofuran	2.093	1.994	4.7	79	0.00
55	P	4-Nitrophenol	0.257	0.260	-1.2	79	0.00
56		2,4-Dinitrotoluene	0.417	0.482	-15.6	90	0.00
57		Fluorene	1.700	1.624	4.5	80	-0.01
58		2,3,4,6-Tetrachlorophenol	0.525	0.527	-0.4	82	-0.01
59		Diethylphthalate	1.717	1.635	4.8	78	-0.01
60		4-Chlorophenyl-phenylether	1.040	0.995	4.3	79	-0.01
61		4-Nitroaniline	0.309	0.338	-9.4	83	-0.01
62		Azobenzene	1.101	1.031	6.4	77	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	81	-0.01
64		4,6-Dinitro-2-methylphenol	0.086	0.099	-15.1	92	0.00
65	c	n-Nitrosodiphenylamine	0.664	0.637	4.1	79	0.00
66		4-Bromophenyl-phenylether	0.289	0.282	2.4	80	-0.01
67		Hexachlorobenzene	0.296	0.301	-1.7	83	-0.01
68		Atrazine	0.258	0.253	1.9	80	-0.01
69	C	Pentachlorophenol	0.203	0.203	0.0	78	0.00
70		Phenanthrene	1.132	1.087	4.0	79	0.00
71		Anthracene	1.142	1.101	3.6	80	-0.01
72		Carbazole	1.011	0.960	5.0	78	-0.01
73		Di-n-butylphthalate	1.123	1.089	3.0	80	-0.02
74	C	Fluoranthene	1.389	1.345	3.2	81	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	84	-0.01
76		Benzidine	0.617	0.438	29.0#	58	0.00
77		Pyrene	1.278	1.217	4.8	81	0.00
78	S	Terphenyl-d14	0.875	0.850	2.9	83	-0.01
79		Butylbenzylphthalate	0.429	0.420	2.1	82	-0.02
80		Benzo(a)anthracene	1.272	1.217	4.3	81	-0.01
81		3,3'-Dichlorobenzidine	0.493	0.478	3.0	81	-0.02
82		Chrysene	1.204	1.148	4.7	81	-0.01
83		Bis(2-ethylhexyl)phthalate	0.621	0.607	2.3	82	-0.02
84	c	Di-n-octyl phthalate	1.046	1.013	3.2	80	-0.03
85		Indeno(1,2,3-cd)pyrene	1.544	1.559	-1.0	84	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	85	-0.02
87		Benzo(b)fluoranthene	1.241	1.204	3.0	85	-0.02

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88	Benzo(k)fluoranthene	1.242	1.174	5.5	81	-0.02
89 C	Benzo(a)pyrene	1.189	1.138	4.3	82	-0.02
90	Dibenzo(a,h)anthracene	1.228	1.196	2.6	84	-0.05
91	Benzo(g,h,i)perylene	1.450	1.422	1.9	84	-0.05

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