Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG122217\
Data File : BG031733.D
Acq On : 23 Dec 2017 4:48
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 25 04:39:11 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	84	-0.01
2		1,4-Dioxane	0.407	0.367	9.8	79	0.00
3		Pyridine	1.088	1.090	-0.2	80	-0.01
4		n-Nitrosodimethylamine	0.439	0.416	5.2	79	0.00
5	S	2-Fluorophenol	1.098	1.064	3.1	82	-0.01
6		Aniline	1.834	1.726	5.9	77	0.00
7	S	Phenol-d6	1.426	1.422	0.3	82	0.00
8		2-Chlorophenol	1.274	1.230	3.5	81	-0.01
9		Benzaldehyde	0.858	0.790	7.9	76	-0.01
10	C	Phenol	1.439	1.429	0.7	83	0.00
11		bis(2-Chloroethyl)ether	1.195	1.104	7.6	78	-0.01
12		1,3-Dichlorobenzene	1.582	1.518	4.0	81	-0.01
13	C	1,4-Dichlorobenzene	1.616	1.541	4.6	82	-0.01
14		1,2-Dichlorobenzene	1.531	1.465	4.3	81	-0.01
15		Benzyl Alcohol	1.070	1.060	0.9	80	-0.01
16		2,2'-oxybis(1-Chloropropane	0.973	0.835	14.2	73	0.00
17		2-Methylphenol	1.030	1.006	2.3	81	0.00
18		Hexachloroethane	0.489	0.475	2.9	84	-0.02
19	P	n-Nitroso-di-n-propylamine	0.974	0.894	8.2	77	-0.02
20		3+4-Methylphenols	1.464	1.440	1.6	81	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	82	-0.01
22		Acetophenone	0.465	0.445	4.3	79	-0.01
23	S	Nitrobenzene-d5	0.265	0.278	-4.9	84	-0.01
24		Nitrobenzene	0.273	0.272	0.4	79	-0.01
25		Isophorone	0.570	0.529	7.2	76	-0.01
26	C	2-Nitrophenol	0.143	0.170	-18.9	96	-0.01
27		2,4-Dimethylphenol	0.301	0.294	2.3	82	-0.01
28		bis(2-Chloroethoxy)methane	0.383	0.353	7.8	76	-0.01
29	C	2,4-Dichlorophenol	0.354	0.357	-0.8	82	0.00
30		1,2,4-Trichlorobenzene	0.432	0.419	3.0	80	-0.01
31		Naphthalene	1.009	0.970	3.9	79	-0.01
32		Benzoic acid	0.186	0.209	-12.4	87	-0.02
33		4-Chloroaniline	0.449	0.423	5.8	77	0.00
34	C	Hexachlorobutadiene	0.296	0.289	2.4	81	0.00
35		Caprolactam	0.107	0.106	0.9	79	-0.01
36	C	4-Chloro-3-methylphenol	0.327	0.329	-0.6	83	-0.01
37		2-Methylnaphthalene	0.818	0.787	3.8	80	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	80	0.00
39		1,2,4,5-Tetrachlorobenzene	0.842	0.836	0.7	80	-0.01
40	P	Hexachlorocyclopentadiene	0.444	0.453	-2.0	80	-0.01
41	S	2,4,6-Tribromophenol	0.305	0.339	-11.1	88	-0.01
42	C	2,4,6-Trichlorophenol	0.486	0.509	-4.7	83	0.00
43		2,4,5-Trichlorophenol	0.538	0.569	-5.8	84	-0.01
44	S	2-Fluorobiphenyl	1.593	1.571	1.4	81	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.708	1.700	0.5	80	-0.01
46		2-Chloronaphthalene	1.303	1.296	0.5	81	-0.01
47		2-Nitroaniline	0.225	0.251	-11.6	88	0.00
48		Acenaphthylene	2.084	2.051	1.6	80	-0.01
49		Dimethylphthalate	1.762	1.724	2.2	80	-0.02
50		2,6-Dinitrotoluene	0.324	0.365	-12.7	88	0.00
51	C	Acenaphthene	1.365	1.342	1.7	80	-0.01
52		3-Nitroaniline	0.316	0.345	-9.2	85	0.00
53	P	2,4-Dinitrophenol	0.115	0.107	7.0	76	-0.01
54		Dibenzofuran	2.093	2.060	1.6	80	-0.01
55	P	4-Nitrophenol	0.257	0.266	-3.5	80	-0.01
56		2,4-Dinitrotoluene	0.417	0.495	-18.7	91	0.00
57		Fluorene	1.700	1.676	1.4	81	-0.01
58		2,3,4,6-Tetrachlorophenol	0.525	0.544	-3.6	83	-0.01
59		Diethylphthalate	1.717	1.658	3.4	78	-0.02
60		4-Chlorophenyl-phenylether	1.040	1.017	2.2	80	-0.01
61		4-Nitroaniline	0.309	0.342	-10.7	83	-0.01
62		Azobenzene	1.101	1.048	4.8	77	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	80	-0.01
64		4,6-Dinitro-2-methylphenol	0.086	0.101	-17.4	92	0.00
65	c	n-Nitrosodiphenylamine	0.664	0.645	2.9	78	-0.01
66		4-Bromophenyl-phenylether	0.289	0.291	-0.7	81	-0.01
67		Hexachlorobenzene	0.296	0.308	-4.1	84	-0.01
68		Atrazine	0.258	0.253	1.9	79	-0.01
69	C	Pentachlorophenol	0.203	0.211	-3.9	80	0.00
70		Phenanthrene	1.132	1.108	2.1	80	-0.01
71		Anthracene	1.142	1.103	3.4	79	-0.01
72		Carbazole	1.011	0.966	4.5	77	-0.01
73		Di-n-butylphthalate	1.123	1.092	2.8	78	-0.02
74	C	Fluoranthene	1.389	1.351	2.7	80	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	83	-0.02
76		Benzidine	0.617	0.459	25.6#	61	-0.01
77		Pyrene	1.278	1.211	5.2	80	-0.01
78	S	Terphenyl-d14	0.875	0.853	2.5	83	-0.01
79		Butylbenzylphthalate	0.429	0.424	1.2	82	-0.02
80		Benzo(a)anthracene	1.272	1.229	3.4	82	-0.02
81		3,3'-Dichlorobenzidine	0.493	0.487	1.2	82	-0.02
82		Chrysene	1.204	1.152	4.3	81	-0.02
83		Bis(2-ethylhexyl)phthalate	0.621	0.612	1.4	82	-0.03
84	c	Di-n-octyl phthalate	1.046	1.038	0.8	82	-0.05
85		Indeno(1,2,3-cd)pyrene	1.544	1.594	-3.2	85	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	85	-0.03
87		Benzo(b)fluoranthene	1.241	1.214	2.2	85	-0.03

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88	Benzo(k)fluoranthene	1.242	1.187	4.4	82	-0.03
89 C	Benzo(a)pyrene	1.189	1.153	3.0	83	-0.03
90	Dibenzo(a,h)anthracene	1.228	1.215	1.1	85	-0.05
91	Benzo(g,h,i)perylene	1.450	1.449	0.1	85	-0.05

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