Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG122217\
Data File : BG031752.D
Acq On : 23 Dec 2017 17:04
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 26 00:35:17 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	93	-0.01
2		1,4-Dioxane	0.407	0.353	13.3	84	0.00
3		Pyridine	1.088	1.072	1.5	87	-0.01
4		n-Nitrosodimethylamine	0.439	0.417	5.0	87	0.00
5	S	2-Fluorophenol	1.098	1.072	2.4	91	0.00
6		Aniline	1.834	1.734	5.5	86	0.00
7	S	Phenol-d6	1.426	1.433	-0.5	91	0.00
8		2-Chlorophenol	1.274	1.251	1.8	91	-0.01
9		Benzaldehyde	0.858	0.814	5.1	87	-0.01
10	C	Phenol	1.439	1.406	2.3	90	0.00
11		bis(2-Chloroethyl)ether	1.195	1.089	8.9	84	0.00
12		1,3-Dichlorobenzene	1.582	1.509	4.6	89	-0.01
13	C	1,4-Dichlorobenzene	1.616	1.553	3.9	91	-0.01
14		1,2-Dichlorobenzene	1.531	1.469	4.0	89	0.00
15		Benzyl Alcohol	1.070	1.064	0.6	89	-0.01
16		2,2'-oxybis(1-Chloropropane	0.973	0.831	14.6	80	0.00
17		2-Methylphenol	1.030	1.006	2.3	89	0.00
18		Hexachloroethane	0.489	0.471	3.7	92	-0.02
19	P	n-Nitroso-di-n-propylamine	0.974	0.906	7.0	86	-0.02
20		3+4-Methylphenols	1.464	1.438	1.8	89	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	90	-0.01
22		Acetophenone	0.465	0.446	4.1	87	-0.01
23	S	Nitrobenzene-d5	0.265	0.286	-7.9	94	-0.01
24		Nitrobenzene	0.273	0.284	-4.0	91	-0.01
25		Isophorone	0.570	0.538	5.6	84	-0.01
26	C	2-Nitrophenol	0.143	0.176	-23.1#	109	-0.01
27		2,4-Dimethylphenol	0.301	0.289	4.0	89	-0.01
28		bis(2-Chloroethoxy)methane	0.383	0.354	7.6	84	-0.01
29	C	2,4-Dichlorophenol	0.354	0.361	-2.0	91	0.00
30		1,2,4-Trichlorobenzene	0.432	0.422	2.3	89	-0.01
31		Naphthalene	1.009	0.985	2.4	89	-0.01
32		Benzoic acid	0.186	0.134	28.0#	61	-0.03
33		4-Chloroaniline	0.449	0.430	4.2	86	0.00
34	C	Hexachlorobutadiene	0.296	0.295	0.3	91	-0.01
35		Caprolactam	0.107	0.103	3.7	84	0.00
36	C	4-Chloro-3-methylphenol	0.327	0.325	0.6	90	-0.01
37		2-Methylnaphthalene	0.818	0.796	2.7	89	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
39		1,2,4,5-Tetrachlorobenzene	0.842	0.827	1.8	89	-0.01
40	P	Hexachlorocyclopentadiene	0.444	0.448	-0.9	88	-0.01
41	S	2,4,6-Tribromophenol	0.305	0.334	-9.5	97	-0.01
42	C	2,4,6-Trichlorophenol	0.486	0.502	-3.3	91	0.00
43		2,4,5-Trichlorophenol	0.538	0.562	-4.5	93	-0.01
44	S	2-Fluorobiphenyl	1.593	1.560	2.1	89	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.708	1.682	1.5	88	-0.01
46		2-Chloronaphthalene	1.303	1.269	2.6	88	-0.01
47		2-Nitroaniline	0.225	0.259	-15.1	101	0.00
48		Acenaphthylene	2.084	2.021	3.0	88	-0.01
49		Dimethylphthalate	1.762	1.687	4.3	87	-0.01
50		2,6-Dinitrotoluene	0.324	0.360	-11.1	97	0.00
51	C	Acenaphthene	1.365	1.293	5.3	85	0.00
52		3-Nitroaniline	0.316	0.338	-7.0	93	0.00
53	P	2,4-Dinitrophenol	0.115	0.034#	70.4#	27#	0.00
54		Dibenzofuran	2.093	2.004	4.3	87	-0.01
55	P	4-Nitrophenol	0.257	0.240	6.6	80	0.00
56		2,4-Dinitrotoluene	0.417	0.497	-19.2	102	0.00
57		Fluorene	1.700	1.619	4.8	87	-0.01
58		2,3,4,6-Tetrachlorophenol	0.525	0.522	0.6	89	-0.01
59		Diethylphthalate	1.717	1.636	4.7	86	-0.02
60		4-Chlorophenyl-phenylether	1.040	1.008	3.1	88	-0.01
61		4-Nitroaniline	0.309	0.336	-8.7	91	0.00
62		Azobenzene	1.101	1.019	7.4	83	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	89	-0.01
64		4,6-Dinitro-2-methylphenol	0.086	0.051	40.7#	52	0.00
65	c	n-Nitrosodiphenylamine	0.664	0.632	4.8	86	-0.01
66		4-Bromophenyl-phenylether	0.289	0.287	0.7	89	-0.01
67		Hexachlorobenzene	0.296	0.298	-0.7	91	-0.01
68		Atrazine	0.258	0.244	5.4	85	-0.01
69	C	Pentachlorophenol	0.203	0.160	21.2#	67	0.00
70		Phenanthrene	1.132	1.062	6.2	85	-0.01
71		Anthracene	1.142	1.072	6.1	85	-0.01
72		Carbazole	1.011	0.931	7.9	83	-0.01
73		Di-n-butylphthalate	1.123	1.061	5.5	85	-0.02
74	C	Fluoranthene	1.389	1.296	6.7	85	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	90	-0.01
76		Benzidine	0.617	0.501	18.8	71	-0.01
77		Pyrene	1.278	1.208	5.5	86	-0.01
78	S	Terphenyl-d14	0.875	0.845	3.4	88	-0.01
79		Butylbenzylphthalate	0.429	0.424	1.2	88	-0.02
80		Benzo(a)anthracene	1.272	1.228	3.5	88	-0.01
81		3,3'-Dichlorobenzidine	0.493	0.495	-0.4	90	-0.01
82		Chrysene	1.204	1.156	4.0	87	-0.01
83		Bis(2-ethylhexyl)phthalate	0.621	0.613	1.3	88	-0.03
84	c	Di-n-octyl phthalate	1.046	1.042	0.4	88	-0.04
85		Indeno(1,2,3-cd)pyrene	1.544	1.589	-2.9	91	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	91	-0.02
87		Benzo(b)fluoranthene	1.241	1.215	2.1	92	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.242	1.171	5.7	87	-0.02
89 C	Benzo(a)pyrene	1.189	1.143	3.9	89	-0.02
90	Dibenzo(a,h)anthracene	1.228	1.214	1.1	91	-0.05
91	Benzo(g,h,i)perylene	1.450	1.440	0.7	91	-0.05

(#) = Out of Range SPCC's out = 1 CCC's out = 2