Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG103117\
Data File : BG030611.D
Acq On : 31 Oct 2017 14:19
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 01 06:37:24 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 17:32:15 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	125	-0.01
2		1,4-Dioxane	0.486	0.459	5.6	110	0.00
3		Pyridine	1.415	1.421	-0.4	115	0.00
4		n-Nitrosodimethylamine	0.661	0.625	5.4	110	0.00
5	S	2-Fluorophenol	1.197	1.183	1.2	114	0.00
6		Aniline	2.081	2.076	0.2	114	0.00
7	S	Phenol-d6	1.680	1.686	-0.4	114	0.00
8		2-Chlorophenol	1.372	1.305	4.9	111	-0.01
9		Benzaldehyde	0.845	0.966	-14.3	131	0.00
10	C	Phenol	1.812	1.712	5.5	108	0.00
11		bis(2-Chloroethyl)ether	1.320	1.331	-0.8	114	-0.01
12		1,3-Dichlorobenzene	1.541	1.470	4.6	111	-0.01
13	C	1,4-Dichlorobenzene	1.573	1.508	4.1	111	-0.02
14		1,2-Dichlorobenzene	1.517	1.449	4.5	111	-0.01
15		Benzyl Alcohol	1.184	1.271	-7.3	124	-0.01
16		2,2'-oxybis(1-Chloropropane	2.242	1.831	18.3	94	-0.02
17		2-Methylphenol	1.204	1.185	1.6	113	0.00
18		Hexachloroethane	0.536	0.525	2.1	113	-0.01
19	P	n-Nitroso-di-n-propylamine	1.143	1.208	-5.7	122	-0.01
20		3+4-Methylphenols	1.696	1.704	-0.5	115	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	132	-0.02
22		Acetophenone	0.482	0.461	4.4	119	-0.01
23	S	Nitrobenzene-d5	0.315	0.312	1.0	121	-0.01
24		Nitrobenzene	0.354	0.317	10.5	109	-0.01
25		Isophorone	0.657	0.596	9.3	111	-0.01
26	C	2-Nitrophenol	0.169	0.171	-1.2	120	-0.02
27		2,4-Dimethylphenol	0.306	0.253	17.3	103	-0.01
28		bis(2-Chloroethoxy)methane	0.403	0.391	3.0	120	-0.02
29	C	2,4-Dichlorophenol	0.326	0.308	5.5	117	-0.01
30		1,2,4-Trichlorobenzene	0.353	0.342	3.1	121	-0.01
31		Naphthalene	0.997	0.921	7.6	115	-0.02
32		Benzoic acid	0.256	0.219	14.5	100	-0.01
33		4-Chloroaniline	0.419	0.412	1.7	121	-0.01
34	C	Hexachlorobutadiene	0.219	0.223	-1.8	126	-0.02
35		Caprolactam	0.108	0.105	2.8	121	0.00
36	C	4-Chloro-3-methylphenol	0.359	0.311	13.4	108	-0.01
37		2-Methylnaphthalene	0.726	0.701	3.4	120	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	137	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.730	0.715	2.1	126	-0.02
40	P	Hexachlorocyclopentadiene	0.247	0.288	-16.6	145	-0.02
41	S	2,4,6-Tribromophenol	0.258	0.257	0.4	123	-0.01
42	C	2,4,6-Trichlorophenol	0.458	0.430	6.1	118	-0.01
43		2,4,5-Trichlorophenol	0.471	0.460	2.3	123	-0.01
44	S	2-Fluorobiphenyl	1.364	1.297	4.9	122	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.719	1.592	7.4	118	-0.01
46		2-Chloronaphthalene	1.254	1.143	8.9	116	-0.01
47		2-Nitroaniline	0.344	0.327	4.9	115	-0.01
48		Acenaphthylene	1.962	1.852	5.6	120	-0.01
49		Dimethylphthalate	1.560	1.478	5.3	119	-0.02
50		2,6-Dinitrotoluene	0.327	0.321	1.8	118	-0.01
51	C	Acenaphthene	1.277	1.148	10.1	113	-0.01
52		3-Nitroaniline	0.353	0.337	4.5	117	-0.01
53	P	2,4-Dinitrophenol	0.153	0.123	19.6	101	0.00
54		Dibenzofuran	1.823	1.745	4.3	122	-0.01
55	P	4-Nitrophenol	0.291	0.249	14.4	104	0.00
56		2,4-Dinitrotoluene	0.443	0.439	0.9	120	-0.01
57		Fluorene	1.506	1.369	9.1	116	-0.01
58		2,3,4,6-Tetrachlorophenol	0.393	0.393	0.0	124	-0.01
59		Diethylphthalate	1.555	1.457	6.3	118	-0.01
60		4-Chlorophenyl-phenylether	0.803	0.800	0.4	126	-0.01
61		4-Nitroaniline	0.362	0.335	7.5	116	-0.01
62		Azobenzene	1.311	1.225	6.6	119	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	134	-0.01
64		4,6-Dinitro-2-methylphenol	0.105	0.114	-8.6	132	-0.01
65	c	n-Nitrosodiphenylamine	0.599	0.585	2.3	122	-0.01
66		4-Bromophenyl-phenylether	0.229	0.229	0.0	125	-0.01
67		Hexachlorobenzene	0.260	0.242	6.9	117	-0.01
68		Atrazine	0.214	0.218	-1.9	125	-0.01
69	C	Pentachlorophenol	0.149	0.149	0.0	119	-0.01
70		Phenanthrene	1.033	0.957	7.4	116	-0.01
71		Anthracene	1.037	0.963	7.1	116	-0.02
72		Carbazole	0.997	0.885	11.2	111	0.00
73		Di-n-butylphthalate	1.146	1.067	6.9	115	-0.02
74	C	Fluoranthene	1.208	1.149	4.9	118	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	133	0.00
76		Benzidine	0.604	0.628	-4.0	126	-0.01
77		Pyrene	1.213	1.128	7.0	116	0.00
78	S	Terphenyl-d14	0.879	0.808	8.1	115	-0.01
79		Butylbenzylphthalate	0.517	0.488	5.6	117	-0.01
80		Benzo(a)anthracene	1.174	1.143	2.6	122	-0.01
81		3,3'-Dichlorobenzidine	0.485	0.474	2.3	122	-0.01
82		Chrysene	1.125	1.066	5.2	120	0.00
83		Bis(2-ethylhexyl)phthalate	0.742	0.679	8.5	114	-0.02
84	c	Di-n-octyl phthalate	1.293	1.162	10.1	112	-0.03
85		Indeno(1,2,3-cd)pyrene	1.383	1.361	1.6	123	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	132	-0.01
87		Benzo(b)fluoranthene	1.145	1.128	1.5	121	-0.01

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88	Benzo(k)fluoranthene	1.141	1.092	4.3	119	-0.01
89 C	Benzo(a)pyrene	1.101	1.065	3.3	120	-0.01
90	Dibenzo(a,h)anthracene	1.114	1.117	-0.3	123	0.00
91	Benzo(g,h,i)perylene	1.035	1.061	-2.5	124	0.00

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