Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071116\
Data File : BG023001.D
Acq On : 11 Jul 2016 13:01
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 12 03:48:35 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 07 16:51:07 2016
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	0.00
2		1,4-Dioxane	0.475	0.444	6.5	121	0.00
3		Pyridine	1.627	1.568	3.6	121	0.00
4		n-Nitrosodimethylamine	0.698	0.701	-0.4	124	0.00
5	S	2-Fluorophenol	1.223	1.148	6.1	120	0.00
6		Aniline	2.654	2.664	-0.4	124	0.00
7	S	Phenol-d6	1.915	1.933	-0.9	123	0.00
8		2-Chlorophenol	1.301	1.330	-2.2	128	0.00
9		Benzaldehyde	1.280	1.240	3.1	123	0.00
10	C	Phenol	1.971	2.036	-3.3	128	0.01
11		bis(2-Chloroethyl)ether	1.562	1.521	2.6	122	0.00
12		1,3-Dichlorobenzene	1.593	1.552	2.6	126	0.00
13	C	1,4-Dichlorobenzene	1.595	1.562	2.1	124	0.00
14		1,2-Dichlorobenzene	1.585	1.561	1.5	128	0.00
15		Benzyl Alcohol	1.754	1.814	-3.4	125	0.00
16		2,2'-oxybis(1-Chloropropane	1.908	1.921	-0.7	126	0.01
17		2-Methylphenol	1.283	1.300	-1.3	126	0.00
18		Hexachloroethane	0.673	0.655	2.7	131	0.00
19	P	n-Nitroso-di-n-propylamine	1.727	1.696	1.8	119	0.00
20		3+4-Methylphenols	1.789	1.875	-4.8	123	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	124	0.00
22		Acetophenone	0.600	0.568	5.3	118	0.00
23	S	Nitrobenzene-d5	0.467	0.461	1.3	124	0.00
24		Nitrobenzene	0.496	0.494	0.4	127	0.00
25		Isophorone	0.939	0.884	5.9	114	0.00
26	C	2-Nitrophenol	0.195	0.191	2.1	116	0.00
27		2,4-Dimethylphenol	0.337	0.312	7.4	116	0.00
28		bis(2-Chloroethoxy)methane	0.524	0.495	5.5	116	0.00
29	C	2,4-Dichlorophenol	0.359	0.363	-1.1	123	0.00
30		1,2,4-Trichlorobenzene	0.411	0.400	2.7	122	0.01
31		Naphthalene	1.018	1.018	0.0	126	0.00
32		Benzoic acid	0.306	0.302	1.3	116	0.03
33		4-Chloroaniline	0.498	0.486	2.4	122	0.01
34	C	Hexachlorobutadiene	0.334	0.313	6.3	120	0.00
35		Caprolactam	0.148	0.130	12.2	109	0.02
36	C	4-Chloro-3-methylphenol	0.491	0.480	2.2	119	0.01
37		2-Methylnaphthalene	0.805	0.809	-0.5	123	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	118	0.00
39		1,2,4,5-Tetrachlorobenzene	0.862	0.855	0.8	119	0.00
40	P	Hexachlorocyclopentadiene	0.496	0.389	21.6	93	0.00
41	S	2,4,6-Tribromophenol	0.359	0.298	17.0	101	0.00
42	C	2,4,6-Trichlorophenol	0.495	0.490	1.0	116	0.00
43		2,4,5-Trichlorophenol	0.509	0.488	4.1	116	0.00
44	S	2-Fluorobiphenyl	1.270	1.202	5.4	115	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.501	1.487	0.9	118	0.00
46		2-Chloronaphthalene	1.217	1.215	0.2	119	0.00
47		2-Nitroaniline	0.519	0.513	1.2	118	0.01
48		Acenaphthylene	1.826	1.763	3.5	115	0.00
49		Dimethylphthalate	1.634	1.470	10.0	109	0.00
50		2,6-Dinitrotoluene	0.367	0.348	5.2	113	0.00
51	C	Acenaphthene	1.193	1.156	3.1	117	0.00
52		3-Nitroaniline	0.366	0.335	8.5	108	0.00
53	P	2,4-Dinitrophenol	0.252	0.192	23.8	87	0.00
54		Dibenzofuran	1.786	1.650	7.6	112	0.00
55	P	4-Nitrophenol	0.307	0.269	12.4	104	0.01
56		2,4-Dinitrotoluene	0.535	0.485	9.3	109	0.00
57		Fluorene	1.537	1.415	7.9	112	0.00
58		2,3,4,6-Tetrachlorophenol	0.509	0.450	11.6	105	0.00
59		Diethylphthalate	1.662	1.472	11.4	107	0.00
60		4-Chlorophenyl-phenylether	0.936	0.844	9.8	107	0.00
61		4-Nitroaniline	0.396	0.357	9.8	109	0.00
62		Azobenzene	1.590	1.524	4.2	116	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
64		4,6-Dinitro-2-methylphenol	0.149	0.130	12.8	88	0.00
65	c	n-Nitrosodiphenylamine	0.576	0.600	-4.2	109	0.00
66		4-Bromophenyl-phenylether	0.260	0.259	0.4	105	0.00
67		Hexachlorobenzene	0.283	0.283	0.0	105	0.00
68		Atrazine	0.256	0.246	3.9	103	0.00
69	C	Pentachlorophenol	0.183	0.172	6.0	94	0.00
70		Phenanthrene	0.980	0.958	2.2	105	0.00
71		Anthracene	0.992	0.974	1.8	106	0.00
72		Carbazole	0.858	0.840	2.1	106	0.00
73		Di-n-butylphthalate	0.954	0.980	-2.7	107	0.00
74	C	Fluoranthene	1.203	1.127	6.3	105	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	105	0.02
76		Benzidine	0.573	0.603	-5.2	107	0.00
77		Pyrene	1.124	1.057	6.0	105	0.00
78	S	Terphenyl-d14	0.646	0.626	3.1	102	0.00
79		Butylbenzylphthalate	0.445	0.457	-2.7	109	0.01
80		Benzo(a)anthracene	1.119	1.103	1.4	106	0.02
81		3,3'-Dichlorobenzidine	0.491	0.474	3.5	100	0.02
82		Chrysene	1.050	1.045	0.5	108	0.02
83		Bis(2-ethylhexyl)phthalate	0.618	0.627	-1.5	109	0.02
84	c	Di-n-octyl phthalate	1.031	1.042	-1.1	107	0.03
85		Indeno(1,2,3-cd)pyrene	1.338	1.300	2.8	105	0.11
86	I	Perylene-d12	1.000	1.000	0.0	103	0.06
87		Benzo(b)fluoranthene	1.154	1.172	-1.6	107	0.04

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88	Benzo(k)fluoranthene	1.095	1.097	-0.2	102	0.05
89 C	Benzo(a)pyrene	1.106	1.123	-1.5	106	0.05
90	Dibenzo(a,h)anthracene	1.098	1.090	0.7	104	0.11
91	Benzo(g,h,i)perylene	1.099	1.092	0.6	103	0.12

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