Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG070816\
Data File : BG022918.D
Acq On : 8 Jul 2016 6:03
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 08 07:12:03 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	100	0.00
2		1,4-Dioxane	0.475	0.479	-0.8	105	0.00
3		Pyridine	1.627	1.680	-3.3	104	0.00
4		n-Nitrosodimethylamine	0.698	0.720	-3.2	102	0.00
5	S	2-Fluorophenol	1.223	1.239	-1.3	104	0.00
6		Aniline	2.654	2.746	-3.5	103	0.00
7	S	Phenol-d6	1.915	1.981	-3.4	102	0.00
8		2-Chlorophenol	1.301	1.347	-3.5	104	0.00
9		Benzaldehyde	1.280	1.265	1.2	101	0.00
10	C	Phenol	1.971	2.035	-3.2	103	0.00
11		bis(2-Chloroethyl)ether	1.562	1.543	1.2	100	0.00
12		1,3-Dichlorobenzene	1.593	1.588	0.3	103	0.00
13	C	1,4-Dichlorobenzene	1.595	1.660	-4.1	106	0.00
14		1,2-Dichlorobenzene	1.585	1.573	0.8	104	0.00
15		Benzyl Alcohol	1.754	1.854	-5.7	103	0.00
16		2,2'-oxybis(1-Chloropropane	1.908	1.917	-0.5	101	0.00
17		2-Methylphenol	1.283	1.298	-1.2	101	0.00
18		Hexachloroethane	0.673	0.668	0.7	107	0.00
19	P	n-Nitroso-di-n-propylamine	1.727	1.728	-0.1	98	0.00
20		3+4-Methylphenols	1.789	1.876	-4.9	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
22		Acetophenone	0.600	0.588	2.0	97	0.00
23	S	Nitrobenzene-d5	0.467	0.464	0.6	99	0.00
24		Nitrobenzene	0.496	0.496	0.0	101	0.00
25		Isophorone	0.939	0.917	2.3	94	0.00
26	C	2-Nitrophenol	0.195	0.194	0.5	93	0.00
27		2,4-Dimethylphenol	0.337	0.333	1.2	98	0.00
28		bis(2-Chloroethoxy)methane	0.524	0.517	1.3	96	0.00
29	C	2,4-Dichlorophenol	0.359	0.368	-2.5	99	0.00
30		1,2,4-Trichlorobenzene	0.411	0.412	-0.2	100	0.00
31		Naphthalene	1.018	1.026	-0.8	101	0.00
32		Benzoic acid	0.306	0.301	1.6	92	0.01
33		4-Chloroaniline	0.498	0.501	-0.6	99	0.00
34	C	Hexachlorobutadiene	0.334	0.338	-1.2	103	0.00
35		Caprolactam	0.148	0.129	12.8	86	0.00
36	C	4-Chloro-3-methylphenol	0.491	0.482	1.8	95	0.00
37		2-Methylnaphthalene	0.805	0.811	-0.7	97	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
39		1,2,4,5-Tetrachlorobenzene	0.862	0.880	-2.1	95	0.00
40	P	Hexachlorocyclopentadiene	0.496	0.508	-2.4	95	0.00
41	S	2,4,6-Tribromophenol	0.359	0.335	6.7	88	0.00
42	C	2,4,6-Trichlorophenol	0.495	0.501	-1.2	92	0.00
43		2,4,5-Trichlorophenol	0.509	0.508	0.2	94	0.00
44	S	2-Fluorobiphenyl	1.270	1.282	-0.9	96	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.501	1.547	-3.1	95	0.00
46		2-Chloronaphthalene	1.217	1.236	-1.6	94	0.00
47		2-Nitroaniline	0.519	0.525	-1.2	94	0.00
48		Acenaphthylene	1.826	1.819	0.4	93	0.00
49		Dimethylphthalate	1.634	1.563	4.3	90	0.00
50		2,6-Dinitrotoluene	0.367	0.355	3.3	90	0.00
51	C	Acenaphthene	1.193	1.355	-13.6	106	0.00
52		3-Nitroaniline	0.366	0.375	-2.5	94	0.00
53	P	2,4-Dinitrophenol	0.252	0.225	10.7	80	0.00
54		Dibenzofuran	1.786	1.755	1.7	93	0.00
55	P	4-Nitrophenol	0.307	0.286	6.8	86	0.00
56		2,4-Dinitrotoluene	0.535	0.520	2.8	91	0.00
57		Fluorene	1.537	1.505	2.1	93	0.00
58		2,3,4,6-Tetrachlorophenol	0.509	0.496	2.6	90	0.00
59		Diethylphthalate	1.662	1.571	5.5	89	0.00
60		4-Chlorophenyl-phenylether	0.936	0.902	3.6	89	0.00
61		4-Nitroaniline	0.396	0.374	5.6	89	0.00
62		Azobenzene	1.590	1.555	2.2	92	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
64		4,6-Dinitro-2-methylphenol	0.149	0.146	2.0	84	0.00
65	c	n-Nitrosodiphenylamine	0.576	0.587	-1.9	90	0.00
66		4-Bromophenyl-phenylether	0.260	0.255	1.9	87	0.00
67		Hexachlorobenzene	0.283	0.286	-1.1	90	0.00
68		Atrazine	0.256	0.252	1.6	89	0.00
69	C	Pentachlorophenol	0.183	0.176	3.8	81	0.00
70		Phenanthrene	0.980	0.981	-0.1	91	0.00
71		Anthracene	0.992	0.998	-0.6	92	0.00
72		Carbazole	0.858	0.854	0.5	91	0.00
73		Di-n-butylphthalate	0.954	0.987	-3.5	91	0.00
74	C	Fluoranthene	1.203	1.151	4.3	90	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
76		Benzidine	0.573	0.519	9.4	79	0.00
77		Pyrene	1.124	1.079	4.0	91	0.00
78	S	Terphenyl-d14	0.646	0.665	-2.9	93	0.00
79		Butylbenzylphthalate	0.445	0.442	0.7	89	0.00
80		Benzo(a)anthracene	1.119	1.118	0.1	92	0.00
81		3,3'-Dichlorobenzidine	0.491	0.488	0.6	88	0.00
82		Chrysene	1.050	1.049	0.1	92	0.00
83		Bis(2-ethylhexyl)phthalate	0.618	0.609	1.5	90	0.00
84	c	Di-n-octyl phthalate	1.031	1.030	0.1	90	0.00
85		Indeno(1,2,3-cd)pyrene	1.338	1.309	2.2	90	0.02
86	I	Perylene-d12	1.000	1.000	0.0	88	0.01
87		Benzo(b)fluoranthene	1.154	1.161	-0.6	91	0.01

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88	Benzo(k)fluoranthene	1.095	1.129	-3.1	90	0.01
89 C	Benzo(a)pyrene	1.106	1.105	0.1	89	0.00
90	Dibenzo(a,h)anthracene	1.098	1.104	-0.5	90	0.03
91	Benzo(g,h,i)perylene	1.099	1.114	-1.4	90	0.03

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