Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022716\
Data File : BG021116.D
Acq On : 27 Feb 2016 15:46
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 29 00:16:33 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 26 16:39:44 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	107	0.00
2		1,4-Dioxane	0.529	0.556	-5.1	109	-0.01
3		Pyridine	1.547	1.605	-3.7	106	-0.01
4		n-Nitrosodimethylamine	0.784	0.784	0.0	100	-0.01
5	S	2-Fluorophenol	1.142	1.188	-4.0	107	0.00
6		Aniline	2.103	2.319	-10.3	110	-0.01
7	S	Phenol-d6	1.700	1.847	-8.6	110	-0.02
8		2-Chlorophenol	1.425	1.475	-3.5	105	-0.01
9		Benzaldehyde	0.921	1.044	-13.4	118	0.00
10	C	Phenol	1.741	1.954	-12.2	113	-0.01
11		bis(2-Chloroethyl)ether	1.335	1.514	-13.4	117	-0.01
12		1,3-Dichlorobenzene	1.516	1.543	-1.8	104	0.00
13	C	1,4-Dichlorobenzene	1.620	1.589	1.9	104	-0.01
14		1,2-Dichlorobenzene	1.534	1.528	0.4	102	-0.01
15		Benzyl Alcohol	1.485	1.605	-8.1	108	-0.01
16		2,2'-oxybis(1-Chloropropane	1.530	2.113	-38.1#	144	0.00
17		2-Methylphenol	1.229	1.328	-8.1	108	-0.01
18		Hexachloroethane	0.613	0.636	-3.8	105	-0.02
19	P	n-Nitroso-di-n-propylamine	1.311	1.529	-16.6	114	-0.02
20		3+4-Methylphenols	1.704	1.822	-6.9	107	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	105	-0.01
22		Acetophenone	0.552	0.573	-3.8	106	-0.01
23	S	Nitrobenzene-d5	0.420	0.461	-9.8	112	-0.02
24		Nitrobenzene	0.434	0.487	-12.2	113	-0.01
25		Isophorone	0.787	0.873	-10.9	113	-0.01
26	C	2-Nitrophenol	0.183	0.183	0.0	102	-0.01
27		2,4-Dimethylphenol	0.323	0.318	1.5	100	-0.02
28		bis(2-Chloroethoxy)methane	0.383	0.425	-11.0	115	-0.01
29	C	2,4-Dichlorophenol	0.317	0.296	6.6	99	-0.01
30		1,2,4-Trichlorobenzene	0.359	0.324	9.7	92	-0.01
31		Naphthalene	1.027	1.036	-0.9	104	-0.01
32		Benzoic acid	0.297	0.274	7.7	91	-0.04
33		4-Chloroaniline	0.429	0.444	-3.5	106	-0.01
34	C	Hexachlorobutadiene	0.240	0.222	7.5	91	-0.01
35		Caprolactam	0.107	0.121	-13.1	114	-0.03
36	C	4-Chloro-3-methylphenol	0.388	0.407	-4.9	106	0.00
37		2-Methylnaphthalene	0.721	0.709	1.7	101	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
39		1,2,4,5-Tetrachlorobenzene	0.739	0.645	12.7	91	-0.01
40	P	Hexachlorocyclopentadiene	0.486	0.411	15.4	87	-0.02
41	S	2,4,6-Tribromophenol	0.262	0.227	13.4	88	-0.01
42	C	2,4,6-Trichlorophenol	0.463	0.443	4.3	98	0.00
43		2,4,5-Trichlorophenol	0.476	0.471	1.1	99	-0.01
44	S	2-Fluorobiphenyl	1.563	1.423	9.0	97	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.684	1.611	4.3	102	-0.01
46		2-Chloronaphthalene	1.278	1.224	4.2	102	-0.01
47		2-Nitroaniline	0.414	0.504	-21.7	125	0.00
48		Acenaphthylene	2.022	2.001	1.0	106	-0.01
49		Dimethylphthalate	1.724	1.687	2.1	104	-0.01
50		2,6-Dinitrotoluene	0.355	0.358	-0.8	105	-0.01
51	C	Acenaphthene	1.379	1.350	2.1	102	0.00
52		3-Nitroaniline	0.345	0.380	-10.1	115	-0.01
53	P	2,4-Dinitrophenol	0.247	0.243	1.6	99	0.00
54		Dibenzofuran	1.747	1.691	3.2	102	-0.01
55	P	4-Nitrophenol	0.306	0.332	-8.5	115	-0.01
56		2,4-Dinitrotoluene	0.509	0.525	-3.1	106	-0.01
57		Fluorene	1.700	1.647	3.1	103	0.00
58		2,3,4,6-Tetrachlorophenol	0.414	0.394	4.8	98	-0.01
59		Diethylphthalate	1.817	1.849	-1.8	109	-0.01
60		4-Chlorophenyl-phenylether	0.935	0.855	8.6	97	-0.01
61		4-Nitroaniline	0.393	0.428	-8.9	116	0.00
62		Azobenzene	1.574	1.840	-16.9	124	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	-0.01
64		4,6-Dinitro-2-methylphenol	0.147	0.143	2.7	102	-0.01
65	c	n-Nitrosodiphenylamine	0.590	0.592	-0.3	107	-0.01
66		4-Bromophenyl-phenylether	0.235	0.212	9.8	95	0.00
67		Hexachlorobenzene	0.248	0.214	13.7	92	0.00
68		Atrazine	0.219	0.221	-0.9	105	-0.01
69	C	Pentachlorophenol	0.168	0.151	10.1	94	-0.01
70		Phenanthrene	1.076	1.074	0.2	107	0.00
71		Anthracene	1.084	1.094	-0.9	107	-0.01
72		Carbazole	0.936	0.990	-5.8	112	0.00
73		Di-n-butylphthalate	1.233	1.345	-9.1	114	-0.02
74	C	Fluoranthene	1.350	1.323	2.0	105	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	100	-0.02
76		Benzidine	0.526	0.578	-9.9	104	0.00
77		Pyrene	1.107	1.190	-7.5	106	-0.01
78	S	Terphenyl-d14	0.856	0.871	-1.8	99	-0.01
79		Butylbenzylphthalate	0.462	0.539	-16.7	111	-0.01
80		Benzo(a)anthracene	1.151	1.168	-1.5	99	-0.01
81		3,3'-Dichlorobenzidine	0.450	0.455	-1.1	95	-0.01
82		Chrysene	1.108	1.100	0.7	98	-0.01
83		Bis(2-ethylhexyl)phthalate	0.714	0.798	-11.8	106	-0.02
84	c	Di-n-octyl phthalate	1.196	1.328	-11.0	105	-0.02
85		Indeno(1,2,3-cd)pyrene	1.336	1.263	5.5	92	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	95	-0.02
87		Benzo(b)fluoranthene	1.250	1.268	-1.4	94	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.232	1.266	-2.8	98	-0.02
89 C	Benzo(a)pyrene	1.200	1.205	-0.4	94	-0.02
90	Dibenzo(a,h)anthracene	1.210	1.192	1.5	92	-0.05
91	Benzo(g,h,i)perylene	1.165	1.142	2.0	92	-0.03

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