Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG093016\
Data File : BG024158.D
Acq On : 30 Sep 2016 14:00
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 30 17:15:20 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG092316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 23 15:35:28 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	108	0.00
2		1,4-Dioxane	0.442	0.477	-7.9	105	-0.01
3		Pyridine	1.562	1.526	2.3	99	0.00
4		n-Nitrosodimethylamine	0.792	0.682	13.9	92	-0.01
5	S	2-Fluorophenol	1.154	1.109	3.9	96	0.00
6		Aniline	2.728	2.419	11.3	96	0.00
7	S	Phenol-d6	1.942	1.827	5.9	100	0.00
8		2-Chlorophenol	1.382	1.237	10.5	97	-0.01
9		Benzaldehyde	1.058	1.082	-2.3	110	0.00
10	C	Phenol	2.042	1.907	6.6	100	0.00
11		bis(2-Chloroethyl)ether	1.610	1.494	7.2	98	0.00
12		1,3-Dichlorobenzene	1.541	1.399	9.2	97	0.00
13	C	1,4-Dichlorobenzene	1.573	1.478	6.0	101	0.00
14		1,2-Dichlorobenzene	1.531	1.440	5.9	99	0.00
15		Benzyl Alcohol	1.880	1.688	10.2	99	0.00
16		2,2'-oxybis(1-Chloropropane	2.223	2.062	7.2	99	-0.01
17		2-Methylphenol	1.358	1.311	3.5	104	0.00
18		Hexachloroethane	0.663	0.604	8.9	96	0.00
19	P	n-Nitroso-di-n-propylamine	1.908	1.676	12.2	96	-0.01
20		3+4-Methylphenols	1.972	1.834	7.0	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	-0.01
22		Acetophenone	0.577	0.540	6.4	101	0.00
23	S	Nitrobenzene-d5	0.507	0.459	9.5	97	0.00
24		Nitrobenzene	0.544	0.481	11.6	95	0.00
25		Isophorone	1.035	0.930	10.1	99	-0.01
26	C	2-Nitrophenol	0.193	0.177	8.3	98	0.00
27		2,4-Dimethylphenol	0.323	0.293	9.3	97	0.00
28		bis(2-Chloroethoxy)methane	0.503	0.477	5.2	103	0.00
29	C	2,4-Dichlorophenol	0.336	0.320	4.8	99	0.00
30		1,2,4-Trichlorobenzene	0.359	0.323	10.0	96	0.00
31		Naphthalene	1.013	0.949	6.3	102	-0.01
32		Benzoic acid	0.241	0.213	11.6	93	-0.01
33		4-Chloroaniline	0.452	0.424	6.2	102	0.00
34	C	Hexachlorobutadiene	0.283	0.243	14.1	93	-0.01
35		Caprolactam	0.137	0.121	11.7	101	-0.01
36	C	4-Chloro-3-methylphenol	0.472	0.436	7.6	97	-0.01
37		2-Methylnaphthalene	0.776	0.718	7.5	100	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	110	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.584	0.534	8.6	96	-0.01
40	P	Hexachlorocyclopentadiene	0.220	0.208	5.5	92	-0.01
41	S	2,4,6-Tribromophenol	0.287	0.232	19.2	88	-0.01
42	C	2,4,6-Trichlorophenol	0.378	0.344	9.0	94	0.00
43		2,4,5-Trichlorophenol	0.389	0.352	9.5	94	0.00
44	S	2-Fluorobiphenyl	1.190	1.114	6.4	99	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.368	1.322	3.4	102	0.00
46		2-Chloronaphthalene	1.021	0.957	6.3	100	-0.01
47		2-Nitroaniline	0.475	0.410	13.7	91	0.00
48		Acenaphthylene	1.734	1.647	5.0	103	-0.01
49		Dimethylphthalate	1.534	1.416	7.7	100	-0.01
50		2,6-Dinitrotoluene	0.322	0.317	1.6	104	-0.01
51	C	Acenaphthene	1.051	0.992	5.6	101	-0.02
52		3-Nitroaniline	0.322	0.314	2.5	107	-0.01
53	P	2,4-Dinitrophenol	0.190	0.161	15.3	89	0.00
54		Dibenzofuran	1.637	1.535	6.2	101	-0.01
55	P	4-Nitrophenol	0.225	0.199	11.6	95	0.00
56		2,4-Dinitrotoluene	0.490	0.452	7.8	101	-0.01
57		Fluorene	1.425	1.319	7.4	101	-0.01
58		2,3,4,6-Tetrachlorophenol	0.375	0.334	10.9	92	-0.01
59		Diethylphthalate	1.652	1.484	10.2	101	0.00
60		4-Chlorophenyl-phenylether	0.763	0.662	13.2	96	-0.01
61		4-Nitroaniline	0.333	0.294	11.7	98	-0.01
62		Azobenzene	1.679	1.521	9.4	99	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	105	-0.01
64		4,6-Dinitro-2-methylphenol	0.139	0.129	7.2	92	0.00
65	c	n-Nitrosodiphenylamine	0.547	0.556	-1.6	101	-0.01
66		4-Bromophenyl-phenylether	0.232	0.212	8.6	93	-0.01
67		Hexachlorobenzene	0.261	0.234	10.3	91	0.00
68		Atrazine	0.242	0.209	13.6	88	-0.01
69	C	Pentachlorophenol	0.124	0.105	15.3	86	0.00
70		Phenanthrene	1.019	0.977	4.1	98	-0.01
71		Anthracene	1.046	0.991	5.3	98	0.00
72		Carbazole	0.974	0.928	4.7	99	-0.01
73		Di-n-butylphthalate	1.275	1.134	11.1	94	-0.01
74	C	Fluoranthene	1.330	1.169	12.1	96	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	131	-0.02
76		Benzidine	0.566	0.430	24.0	91	0.00
77		Pyrene	1.397	1.124	19.5	96	-0.01
78	S	Terphenyl-d14	0.977	0.745	23.7	93	-0.01
79		Butylbenzylphthalate	0.514	0.472	8.2	116	-0.01
80		Benzo(a)anthracene	1.139	1.061	6.8	118	-0.02
81		3,3'-Dichlorobenzidine	0.408	0.397	2.7	121	-0.02
82		Chrysene	1.070	1.008	5.8	121	-0.02
83		Bis(2-ethylhexyl)phthalate	0.614	0.664	-8.1	132	-0.02
84	c	Di-n-octyl phthalate	1.060	1.145	-8.0	128	-0.02
85		Indeno(1,2,3-cd)pyrene	1.193	1.253	-5.0	125	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	138	-0.04
87		Benzo(b)fluoranthene	1.164	1.086	6.7	124	-0.03

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88	Benzo(k)fluoranthene	1.156	1.097	5.1	126	-0.03
89 C	Benzo(a)pyrene	1.125	1.057	6.0	126	-0.03
90	Dibenzo(a,h)anthracene	1.123	1.027	8.5	124	-0.06
91	Benzo(g,h,i)perylene	1.132	1.054	6.9	126	-0.06

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