Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG040715\
Data File : BG016256.D
Acq On : 7 Apr 2015 14:50
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 08 02:33:39 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 07 04:17:17 2015
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	113	0.00
2		1,4-Dioxane	0.576	0.514	10.8	104	0.00
3		Pyridine	1.722	1.657	3.8	109	0.00
4		n-Nitrosodimethylamine	0.512	0.470	8.2	103	0.00
5	S	2-Fluorophenol	1.267	1.223	3.5	110	0.00
6		Aniline	2.717	2.615	3.8	108	0.00
7	S	Phenol-d6	1.902	1.890	0.6	112	0.00
8		2-Chlorophenol	1.522	1.518	0.3	113	0.00
9		Benzaldehyde	1.114	0.997	10.5	106	0.00
10	C	Phenol	2.035	2.014	1.0	111	0.00
11		bis(2-Chloroethyl)ether	1.623	1.548	4.6	110	0.00
12		1,3-Dichlorobenzene	1.537	1.493	2.9	112	0.00
13	C	1,4-Dichlorobenzene	1.594	1.528	4.1	113	0.00
14		1,2-Dichlorobenzene	1.525	1.477	3.1	112	0.00
15		Benzyl Alcohol	1.326	1.348	-1.7	111	0.00
16		2,2'-oxybis(1-Chloropropane	1.828	1.716	6.1	107	0.00
17		2-Methylphenol	1.365	1.362	0.2	111	0.00
18		Hexachloroethane	0.610	0.575	5.7	110	0.00
19	P	n-Nitroso-di-n-propylamine	1.363	1.338	1.8	107	0.00
20		3+4-Methylphenols	1.891	1.947	-3.0	113	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.464	0.438	5.6	110	0.00
23	S	Nitrobenzene-d5	0.345	0.328	4.9	109	0.00
24		Nitrobenzene	0.369	0.345	6.5	108	0.00
25		Isophorone	0.767	0.727	5.2	109	0.00
26	C	2-Nitrophenol	0.172	0.186	-8.1	118	0.00
27		2,4-Dimethylphenol	0.321	0.314	2.2	112	0.00
28		bis(2-Chloroethoxy)methane	0.438	0.413	5.7	110	0.00
29	C	2,4-Dichlorophenol	0.254	0.260	-2.4	115	0.00
30		1,2,4-Trichlorobenzene	0.265	0.257	3.0	114	0.00
31		Naphthalene	1.042	0.986	5.4	111	0.00
32		Benzoic acid	0.182	0.216	-18.7	132	0.00
33		4-Chloroaniline	0.489	0.483	1.2	113	0.00
34	C	Hexachlorobutadiene	0.116	0.116	0.0	118	0.00
35		Caprolactam	0.160	0.165	-3.1	111	0.00
36	C	4-Chloro-3-methylphenol	0.335	0.336	-0.3	113	0.00
37		2-Methylnaphthalene	0.750	0.722	3.7	113	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	112	0.00
39		1,2,4,5-Tetrachlorobenzene	0.440	0.428	2.7	114	0.00
40	P	Hexachlorocyclopentadiene	0.202	0.200	1.0	114	0.00
41	S	2,4,6-Tribromophenol	0.135	0.140	-3.7	115	0.00
42	C	2,4,6-Trichlorophenol	0.327	0.338	-3.4	116	0.00
43		2,4,5-Trichlorophenol	0.350	0.364	-4.0	119	0.00
44	S	2-Fluorobiphenyl	1.175	1.123	4.4	113	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.534	1.463	4.6	113	0.00
46		2-Chloronaphthalene	1.169	1.104	5.6	113	0.00
47		2-Nitroaniline	0.377	0.374	0.8	110	0.00
48		Acenaphthylene	2.078	1.984	4.5	112	0.00
49		Dimethylphthalate	1.466	1.382	5.7	109	0.00
50		2,6-Dinitrotoluene	0.355	0.349	1.7	112	0.00
51	C	Acenaphthene	1.242	1.169	5.9	112	0.00
52		3-Nitroaniline	0.453	0.442	2.4	110	0.00
53	P	2,4-Dinitrophenol	0.167	0.163	2.4	109	0.00
54		Dibenzofuran	1.723	1.637	5.0	113	0.00
55	P	4-Nitrophenol	0.374	0.364	2.7	110	0.00
56		2,4-Dinitrotoluene	0.483	0.483	0.0	112	0.00
57		Fluorene	1.520	1.434	5.7	111	0.00
58		2,3,4,6-Tetrachlorophenol	0.270	0.282	-4.4	120	0.00
59		Diethylphthalate	1.560	1.467	6.0	109	0.00
60		4-Chlorophenyl-phenylether	0.622	0.592	4.8	112	0.00
61		4-Nitroaniline	0.514	0.489	4.9	107	0.00
62		Azobenzene	1.661	1.491	10.2	106	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
64		4,6-Dinitro-2-methylphenol	0.119	0.130	-9.2	111	0.00
65	c	n-Nitrosodiphenylamine	0.676	0.661	2.2	110	0.00
66		4-Bromophenyl-phenylether	0.160	0.158	1.3	111	0.00
67		Hexachlorobenzene	0.166	0.164	1.2	112	0.00
68		Atrazine	0.188	0.186	1.1	110	0.00
69	C	Pentachlorophenol	0.101	0.104	-3.0	113	0.00
70		Phenanthrene	1.125	1.064	5.4	109	0.00
71		Anthracene	1.115	1.069	4.1	109	0.00
72		Carbazole	1.119	1.058	5.5	108	0.00
73		Di-n-butylphthalate	1.369	1.306	4.6	106	0.00
74	C	Fluoranthene	1.159	1.090	6.0	107	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	109	0.00
76		Benzidine	0.848	0.751	11.4	96	0.00
77		Pyrene	1.393	1.329	4.6	107	0.00
78	S	Terphenyl-d14	0.855	0.817	4.4	108	0.00
79		Butylbenzylphthalate	0.684	0.701	-2.5	109	0.00
80		Benzo(a)anthracene	1.182	1.133	4.1	109	0.00
81		3,3'-Dichlorobenzidine	0.389	0.389	0.0	110	0.00
82		Chrysene	1.141	1.066	6.6	107	0.00
83		Bis(2-ethylhexyl)phthalate	0.926	0.926	0.0	108	0.00
84	c	Di-n-octyl phthalate	1.509	1.536	-1.8	106	0.00
85		Indeno(1,2,3-cd)pyrene	1.186	1.176	0.8	111	0.00
86	I	Perylene-d12	1.000	1.000	0.0	110	0.00
87		Benzo(b)fluoranthene	1.171	1.094	6.6	107	0.00

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88	Benzo(k)fluoranthene	1.146	1.104	3.7	113	0.00
89 C	Benzo(a)pyrene	1.098	1.050	4.4	109	0.00
90	Dibenzo(a,h)anthracene	1.068	1.020	4.5	110	0.00
91	Benzo(g,h,i)perylene	1.067	1.025	3.9	111	0.00

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