Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG050815\
Data File : BG016970.D
Acq On : 9 May 2015 00:42
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 09 01:36:17 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 08 23:46:35 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	101	-0.01
2		1,4-Dioxane	0.481	0.419	12.9	93	-0.01
3		Pyridine	1.488	1.353	9.1	93	-0.01
4		n-Nitrosodimethylamine	0.888	0.877	1.2	107	-0.01
5	S	2-Fluorophenol	1.149	1.101	4.2	101	-0.01
6		Aniline	2.304	2.238	2.9	102	-0.01
7	S	Phenol-d6	1.641	1.659	-1.1	107	0.00
8		2-Chlorophenol	1.331	1.329	0.2	106	-0.01
9		Benzaldehyde	1.144	1.106	3.3	99	-0.01
10	C	Phenol	1.760	1.785	-1.4	108	0.00
11		bis(2-Chloroethyl)ether	1.361	1.313	3.5	101	-0.02
12		1,3-Dichlorobenzene	1.467	1.354	7.7	99	-0.01
13	C	1,4-Dichlorobenzene	1.486	1.387	6.7	98	-0.01
14		1,2-Dichlorobenzene	1.426	1.345	5.7	100	-0.02
15		Benzyl Alcohol	1.499	1.535	-2.4	108	-0.01
16		2,2'-oxybis(1-Chloropropane	2.519	2.259	10.3	97	-0.02
17		2-Methylphenol	1.239	1.250	-0.9	106	0.00
18		Hexachloroethane	0.611	0.577	5.6	100	-0.02
19	P	n-Nitroso-di-n-propylamine	1.439	1.398	2.8	104	-0.01
20		3+4-Methylphenols	1.708	1.737	-1.7	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	-0.01
22		Acetophenone	0.476	0.458	3.8	108	-0.02
23	S	Nitrobenzene-d5	0.409	0.402	1.7	111	-0.01
24		Nitrobenzene	0.413	0.400	3.1	109	-0.01
25		Isophorone	0.827	0.778	5.9	107	-0.01
26	C	2-Nitrophenol	0.168	0.173	-3.0	115	-0.01
27		2,4-Dimethylphenol	0.305	0.294	3.6	106	-0.01
28		bis(2-Chloroethoxy)methane	0.417	0.387	7.2	107	-0.02
29	C	2,4-Dichlorophenol	0.288	0.284	1.4	111	-0.01
30		1,2,4-Trichlorobenzene	0.305	0.283	7.2	104	-0.01
31		Naphthalene	0.995	0.923	7.2	106	-0.01
32		Benzoic acid	0.179	0.226	-26.3#	144	0.01
33		4-Chloroaniline	0.459	0.449	2.2	109	-0.01
34	C	Hexachlorobutadiene	0.191	0.168	12.0	101	-0.02
35		Caprolactam	0.150	0.146	2.7	110	0.00
36	C	4-Chloro-3-methylphenol	0.369	0.369	0.0	113	0.00
37		2-Methylnaphthalene	0.758	0.716	5.5	108	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	113	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.541	0.491	9.2	107	-0.01
40	P	Hexachlorocyclopentadiene	0.349	0.294	15.8	100	-0.01
41	S	2,4,6-Tribromophenol	0.201	0.209	-4.0	124	-0.01
42	C	2,4,6-Trichlorophenol	0.383	0.372	2.9	109	-0.01
43		2,4,5-Trichlorophenol	0.407	0.397	2.5	113	0.00
44	S	2-Fluorobiphenyl	1.325	1.186	10.5	106	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.432	1.319	7.9	109	-0.01
46		2-Chloronaphthalene	1.134	1.049	7.5	109	-0.01
47		2-Nitroaniline	0.395	0.442	-11.9	129	0.00
48		Acenaphthylene	1.975	1.828	7.4	109	-0.02
49		Dimethylphthalate	1.493	1.393	6.7	111	-0.01
50		2,6-Dinitrotoluene	0.310	0.318	-2.6	117	-0.01
51	C	Acenaphthene	1.175	1.067	9.2	108	-0.01
52		3-Nitroaniline	0.353	0.381	-7.9	121	0.00
53	P	2,4-Dinitrophenol	0.142	0.131	7.7	115	0.00
54		Dibenzofuran	1.669	1.558	6.7	109	-0.01
55	P	4-Nitrophenol	0.299	0.306	-2.3	119	0.00
56		2,4-Dinitrotoluene	0.397	0.445	-12.1	129	0.00
57		Fluorene	1.478	1.397	5.5	111	-0.01
58		2,3,4,6-Tetrachlorophenol	0.351	0.340	3.1	110	0.00
59		Diethylphthalate	1.577	1.478	6.3	111	-0.02
60		4-Chlorophenyl-phenylether	0.742	0.679	8.5	110	-0.02
61		4-Nitroaniline	0.409	0.412	-0.7	117	0.00
62		Azobenzene	1.633	1.553	4.9	113	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	112	-0.02
64		4,6-Dinitro-2-methylphenol	0.099	0.107	-8.1	129	-0.01
65	c	n-Nitrosodiphenylamine	0.611	0.576	5.7	113	-0.01
66		4-Bromophenyl-phenylether	0.201	0.190	5.5	113	-0.02
67		Hexachlorobenzene	0.212	0.206	2.8	117	-0.01
68		Atrazine	0.220	0.199	9.5	106	-0.01
69	C	Pentachlorophenol	0.137	0.130	5.1	109	0.00
70		Phenanthrene	1.029	0.950	7.7	110	-0.01
71		Anthracene	1.038	0.959	7.6	110	-0.01
72		Carbazole	0.987	0.910	7.8	108	-0.01
73		Di-n-butylphthalate	1.280	1.189	7.1	109	-0.03
74	C	Fluoranthene	1.194	1.087	9.0	106	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	103	-0.02
76		Benzidine	0.680	0.580	14.7	89	-0.02
77		Pyrene	1.180	1.135	3.8	107	-0.02
78	S	Terphenyl-d14	0.853	0.836	2.0	107	-0.02
79		Butylbenzylphthalate	0.561	0.547	2.5	107	-0.03
80		Benzo(a)anthracene	1.073	1.035	3.5	106	-0.01
81		3,3'-Dichlorobenzidine	0.419	0.393	6.2	102	-0.02
82		Chrysene	1.005	0.978	2.7	106	-0.02
83		Bis(2-ethylhexyl)phthalate	0.767	0.734	4.3	103	-0.03
84	c	Di-n-octyl phthalate	1.289	1.239	3.9	104	-0.04
85		Indeno(1,2,3-cd)pyrene	1.198	1.179	1.6	112	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	106	-0.03
87		Benzo(b)fluoranthene	1.114	1.035	7.1	105	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.072	1.028	4.1	109	-0.03
89 C	Benzo(a)pyrene	1.039	0.990	4.7	107	-0.03
90	Dibenzo(a,h)anthracene	1.061	1.011	4.7	110	-0.05
91	Benzo(g,h,i)perylene	1.011	0.954	5.6	110	-0.04

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