Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG120715\
Data File : BG019979.D
Acq On : 7 Dec 2015 12:11
Operator : UM/NP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 08 00:04:38 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 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	-0.02
2		1,4-Dioxane	0.431	0.475	-10.2	112	0.00
3		Pyridine	1.309	1.414	-8.0	108	-0.02
4		n-Nitrosodimethylamine	0.689	0.662	3.9	92	-0.01
5	S	2-Fluorophenol	1.107	1.212	-9.5	111	-0.01
6		Aniline	2.164	2.241	-3.6	103	-0.02
7	S	Phenol-d6	1.671	1.797	-7.5	108	-0.01
8		2-Chlorophenol	1.349	1.423	-5.5	106	-0.02
9		Benzaldehyde	1.114	1.039	6.7	95	-0.02
10	C	Phenol	1.758	1.903	-8.2	107	0.00
11		bis(2-Chloroethyl)ether	1.302	1.387	-6.5	108	-0.02
12		1,3-Dichlorobenzene	1.529	1.582	-3.5	105	-0.02
13	C	1,4-Dichlorobenzene	1.580	1.637	-3.6	106	-0.02
14		1,2-Dichlorobenzene	1.507	1.533	-1.7	103	-0.02
15		Benzyl Alcohol	1.466	1.424	2.9	99	-0.02
16		2,2'-oxybis(1-Chloropropane	2.124	2.187	-3.0	105	-0.02
17		2-Methylphenol	1.241	1.298	-4.6	104	-0.02
18		Hexachloroethane	0.591	0.608	-2.9	102	-0.02
19	P	n-Nitroso-di-n-propylamine	1.318	1.254	4.9	97	-0.02
20		3+4-Methylphenols	1.748	1.794	-2.6	100	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	97	-0.02
22		Acetophenone	0.548	0.568	-3.6	99	-0.02
23	S	Nitrobenzene-d5	0.392	0.425	-8.4	103	-0.02
24		Nitrobenzene	0.425	0.461	-8.5	102	-0.02
25		Isophorone	0.840	0.837	0.4	96	-0.02
26	C	2-Nitrophenol	0.164	0.199	-21.3#	114	-0.01
27		2,4-Dimethylphenol	0.342	0.353	-3.2	100	-0.02
28		bis(2-Chloroethoxy)methane	0.411	0.438	-6.6	103	-0.02
29	C	2,4-Dichlorophenol	0.349	0.378	-8.3	102	-0.02
30		1,2,4-Trichlorobenzene	0.395	0.411	-4.1	99	-0.02
31		Naphthalene	1.071	1.124	-4.9	100	-0.02
32		Benzoic acid	0.211	0.281	-33.2#	129	-0.01
33		4-Chloroaniline	0.472	0.482	-2.1	98	-0.02
34	C	Hexachlorobutadiene	0.291	0.288	1.0	95	-0.02
35		Caprolactam	0.130	0.123	5.4	94	-0.03
36	C	4-Chloro-3-methylphenol	0.435	0.430	1.1	96	-0.01
37		2-Methylnaphthalene	0.785	0.782	0.4	96	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	89	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.759	0.776	-2.2	92	-0.02
40	P	Hexachlorocyclopentadiene	0.433	0.385	11.1	78	-0.02
41	S	2,4,6-Tribromophenol	0.306	0.295	3.6	87	-0.02
42	C	2,4,6-Trichlorophenol	0.502	0.509	-1.4	93	-0.02
43		2,4,5-Trichlorophenol	0.531	0.539	-1.5	94	-0.02
44	S	2-Fluorobiphenyl	1.465	1.501	-2.5	93	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.641	1.677	-2.2	92	-0.02
46		2-Chloronaphthalene	1.265	1.294	-2.3	92	-0.02
47		2-Nitroaniline	0.400	0.454	-13.5	100	-0.02
48		Acenaphthylene	2.155	2.167	-0.6	91	-0.02
49		Dimethylphthalate	1.799	1.709	5.0	87	-0.02
50		2,6-Dinitrotoluene	0.338	0.374	-10.7	96	-0.02
51	C	Acenaphthene	1.308	1.299	0.7	89	-0.02
52		3-Nitroaniline	0.351	0.393	-12.0	98	-0.02
53	P	2,4-Dinitrophenol	0.143	0.125	12.6	80	-0.02
54		Dibenzofuran	1.945	1.922	1.2	90	-0.02
55	P	4-Nitrophenol	0.306	0.320	-4.6	92	0.00
56		2,4-Dinitrotoluene	0.472	0.533	-12.9	96	-0.01
57		Fluorene	1.741	1.676	3.7	88	-0.02
58		2,3,4,6-Tetrachlorophenol	0.491	0.481	2.0	86	-0.01
59		Diethylphthalate	1.947	1.763	9.5	84	-0.02
60		4-Chlorophenyl-phenylether	1.000	0.947	5.3	86	-0.02
61		4-Nitroaniline	0.395	0.415	-5.1	92	-0.02
62		Azobenzene	1.620	1.536	5.2	86	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	82	-0.02
64		4,6-Dinitro-2-methylphenol	0.103	0.117	-13.6	92	-0.02
65	c	n-Nitrosodiphenylamine	0.579	0.597	-3.1	85	-0.02
66		4-Bromophenyl-phenylether	0.249	0.252	-1.2	83	-0.02
67		Hexachlorobenzene	0.259	0.262	-1.2	84	-0.02
68		Atrazine	0.251	0.242	3.6	78	-0.02
69	C	Pentachlorophenol	0.151	0.171	-13.2	92	-0.02
70		Phenanthrene	1.132	1.161	-2.6	85	-0.02
71		Anthracene	1.153	1.178	-2.2	84	-0.02
72		Carbazole	1.015	1.040	-2.5	84	-0.02
73		Di-n-butylphthalate	1.245	1.284	-3.1	85	-0.02
74	C	Fluoranthene	1.447	1.478	-2.1	84	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	83	-0.03
76		Benzidine	0.657	0.608	7.5	75	-0.02
77		Pyrene	1.177	1.208	-2.6	84	-0.02
78	S	Terphenyl-d14	0.820	0.828	-1.0	82	-0.02
79		Butylbenzylphthalate	0.461	0.490	-6.3	87	-0.03
80		Benzo(a)anthracene	1.224	1.235	-0.9	84	-0.02
81		3,3'-Dichlorobenzidine	0.466	0.444	4.7	78	-0.03
82		Chrysene	1.162	1.175	-1.1	84	-0.03
83		Bis(2-ethylhexyl)phthalate	0.677	0.705	-4.1	87	-0.04
84	c	Di-n-octyl phthalate	1.100	1.194	-8.5	88	-0.06
85		Indeno(1,2,3-cd)pyrene	1.280	1.369	-7.0	89	-0.09
86	I	Perylene-d12	1.000	1.000	0.0	85	-0.05
87		Benzo(b)fluoranthene	1.268	1.274	-0.5	86	-0.05

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88	Benzo(k)fluoranthene	1.255	1.265	-0.8	87	-0.05
89 C	Benzo(a)pyrene	1.203	1.215	-1.0	87	-0.05
90	Dibenzo(a,h)anthracene	1.189	1.223	-2.9	89	-0.08
91	Benzo(g,h,i)perylene	1.167	1.199	-2.7	88	-0.09

(#) = Out of Range SPCC's out = 0 CCC's out = 1