Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG051215\
Data File : BG017085.D
Acq On : 12 May 2015 16:22
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 13 03:35:08 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 13 03:21:45 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	56	-0.10
2		1,4-Dioxane	0.481	0.389	19.1	48#	-0.11
3		Pyridine	1.488	1.249	16.1	48#	-0.11
4		n-Nitrosodimethylamine	0.888	1.005	-13.2	68	-0.11
5	S	2-Fluorophenol	1.149	1.090	5.1	55	-0.08
6		Aniline	2.304	1.986	13.8	50	-0.10
7	S	Phenol-d6	1.641	1.460	11.0	52	-0.05
8		2-Chlorophenol	1.331	1.264	5.0	56	-0.08
9		Benzaldehyde	1.144	0.928	18.9	46#	-0.10
10	C	Phenol	1.760	1.532	13.0	51	-0.05
11		bis(2-Chloroethyl)ether	1.361	1.106	18.7	47#	-0.10
12		1,3-Dichlorobenzene	1.467	1.382	5.8	56	-0.10
13	C	1,4-Dichlorobenzene	1.486	1.440	3.1	57	-0.09
14		1,2-Dichlorobenzene	1.426	1.374	3.6	56	-0.10
15		Benzyl Alcohol	1.499	1.411	5.9	55	-0.08
16		2,2'-oxybis(1-Chloropropane	2.519	1.570	37.7#	37#	-0.09
17		2-Methylphenol	1.239	1.158	6.5	55	-0.06
18		Hexachloroethane	0.611	0.576	5.7	55	-0.10
19	P	n-Nitroso-di-n-propylamine	1.439	1.191	17.2	49#	-0.09
20		3+4-Methylphenols	1.708	1.619	5.2	55	-0.06
21	I	Naphthalene-d8	1.000	1.000	0.0	51	-0.09
22		Acetophenone	0.476	0.475	0.2	55	-0.09
23	S	Nitrobenzene-d5	0.409	0.419	-2.4	57	-0.09
24		Nitrobenzene	0.413	0.415	-0.5	55	-0.08
25		Isophorone	0.827	0.784	5.2	53	-0.09
26	C	2-Nitrophenol	0.168	0.197	-17.3	64	-0.08
27		2,4-Dimethylphenol	0.305	0.326	-6.9	58	-0.07
28		bis(2-Chloroethoxy)methane	0.417	0.366	12.2	50#	-0.09
29	C	2,4-Dichlorophenol	0.288	0.321	-11.5	61	-0.07
30		1,2,4-Trichlorobenzene	0.305	0.345	-13.1	62	-0.09
31		Naphthalene	0.995	0.993	0.2	56	-0.09
32		Benzoic acid	0.179	0.066	63.1#	21#	-0.09
33		4-Chloroaniline	0.459	0.481	-4.8	57	-0.08
34	C	Hexachlorobutadiene	0.191	0.224	-17.3	66	-0.10
35		Caprolactam	0.150	0.157	-4.7	58	-0.08
36	C	4-Chloro-3-methylphenol	0.369	0.383	-3.8	57	-0.05
37		2-Methylnaphthalene	0.758	0.778	-2.6	58	-0.09
38	I	Acenaphthene-d10	1.000	1.000	0.0	57	-0.08
39		1,2,4,5-Tetrachlorobenzene	0.541	0.588	-8.7	65	-0.08
40	P	Hexachlorocyclopentadiene	0.349	0.367	-5.2	63	-0.08
41	S	2,4,6-Tribromophenol	0.201	0.270	-34.3#	80	-0.07
42	C	2,4,6-Trichlorophenol	0.383	0.465	-21.4#	69	-0.07
43		2,4,5-Trichlorophenol	0.407	0.470	-15.5	67	-0.05
44	S	2-Fluorobiphenyl	1.325	1.403	-5.9	63	-0.08

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.432	1.448	-1.1	60	-0.08
46		2-Chloronaphthalene	1.134	1.164	-2.6	61	-0.08
47		2-Nitroaniline	0.395	0.438	-10.9	64	-0.07
48		Acenaphthylene	1.975	1.950	1.3	58	-0.08
49		Dimethylphthalate	1.493	1.640	-9.8	66	-0.08
50		2,6-Dinitrotoluene	0.310	0.362	-16.8	67	-0.07
51	C	Acenaphthene	1.175	1.139	3.1	58	-0.08
52		3-Nitroaniline	0.353	0.389	-10.2	62	-0.07
53	P	2,4-Dinitrophenol	0.142	0.122	14.1	54	-0.07
54		Dibenzofuran	1.669	1.716	-2.8	60	-0.08
55	P	4-Nitrophenol	0.299	0.310	-3.7	61	-0.01
56		2,4-Dinitrotoluene	0.397	0.497	-25.2#	72	-0.07
57		Fluorene	1.478	1.562	-5.7	63	-0.07
58		2,3,4,6-Tetrachlorophenol	0.351	0.439	-25.1#	71	-0.06
59		Diethylphthalate	1.577	1.722	-9.2	65	-0.08
60		4-Chlorophenyl-phenylether	0.742	0.800	-7.8	65	-0.08
61		4-Nitroaniline	0.409	0.429	-4.9	61	-0.06
62		Azobenzene	1.633	1.534	6.1	56	-0.08
63	I	Phenanthrene-d10	1.000	1.000	0.0	59	-0.08
64		4,6-Dinitro-2-methylphenol	0.099	0.118	-19.2	75	-0.07
65	c	n-Nitrosodiphenylamine	0.611	0.628	-2.8	65	-0.07
66		4-Bromophenyl-phenylether	0.201	0.220	-9.5	69	-0.08
67		Hexachlorobenzene	0.212	0.232	-9.4	70	-0.07
68		Atrazine	0.220	0.238	-8.2	67	-0.07
69	C	Pentachlorophenol	0.137	0.118	13.9	52	-0.06
70		Phenanthrene	1.029	1.055	-2.5	64	-0.07
71		Anthracene	1.038	1.066	-2.7	65	-0.07
72		Carbazole	0.987	0.961	2.6	60	-0.07
73		Di-n-butylphthalate	1.280	1.313	-2.6	64	-0.08
74	C	Fluoranthene	1.194	1.255	-5.1	65	-0.07
75	I	Chrysene-d12	1.000	1.000	0.0	60	-0.07
76		Benzidine	0.680	0.567	16.6	51	-0.07
77		Pyrene	1.180	1.178	0.2	65	-0.07
78	S	Terphenyl-d14	0.853	0.983	-15.2	73	-0.07
79		Butylbenzylphthalate	0.561	0.547	2.5	62	-0.07
80		Benzo(a)anthracene	1.073	1.161	-8.2	69	-0.07
81		3,3'-Dichlorobenzidine	0.419	0.464	-10.7	71	-0.07
82		Chrysene	1.005	1.104	-9.9	70	-0.07
83		Bis(2-ethylhexyl)phthalate	0.767	0.795	-3.7	65	-0.08
84	c	Di-n-octyl phthalate	1.289	1.279	0.8	63	-0.09
85		Indeno(1,2,3-cd)pyrene	1.198	1.403	-17.1	78	-0.14
86	I	Perylene-d12	1.000	1.000	0.0	63	-0.11
87		Benzo(b)fluoranthene	1.114	1.163	-4.4	70	-0.10

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.072	1.141	-6.4	72	-0.10
89 C	Benzo(a)pyrene	1.039	1.082	-4.1	70	-0.11
90	Dibenzo(a,h)anthracene	1.061	1.176	-10.8	76	-0.15
91	Benzo(g,h,i)perylene	1.011	1.096	-8.4	76	-0.16

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