Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG060415\
Data File : BG017456.D
Acq On : 4 Jun 2015 23:37
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 05 04:00:56 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 05 03:51:43 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	91	0.00
2		1,4-Dioxane	0.511	0.505	1.2	94	0.00
3		Pyridine	1.365	1.402	-2.7	89	0.00
4		n-Nitrosodimethylamine	0.895	1.006	-12.4	97	0.00
5	S	2-Fluorophenol	1.135	1.197	-5.5	95	0.00
6		Aniline	2.076	2.131	-2.6	92	0.00
7	S	Phenol-d6	1.547	1.540	0.5	88	-0.01
8		2-Chlorophenol	1.342	1.335	0.5	90	-0.01
9		Benzaldehyde	1.050	1.013	3.5	83	0.00
10	C	Phenol	1.589	1.700	-7.0	93	0.00
11		bis(2-Chloroethyl)ether	1.210	1.265	-4.5	93	0.00
12		1,3-Dichlorobenzene	1.501	1.495	0.4	91	0.00
13	C	1,4-Dichlorobenzene	1.573	1.534	2.5	89	0.00
14		1,2-Dichlorobenzene	1.472	1.450	1.5	89	0.00
15		Benzyl Alcohol	1.440	1.466	-1.8	88	0.00
16		2,2'-oxybis(1-Chloropropane	2.040	2.134	-4.6	93	0.00
17		2-Methylphenol	1.205	1.224	-1.6	90	0.00
18		Hexachloroethane	0.627	0.631	-0.6	89	0.00
19	P	n-Nitroso-di-n-propylamine	1.308	1.314	-0.5	90	-0.01
20		3+4-Methylphenols	1.668	1.653	0.9	87	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	92	-0.01
22		Acetophenone	0.490	0.484	1.2	89	0.00
23	S	Nitrobenzene-d5	0.401	0.412	-2.7	91	0.00
24		Nitrobenzene	0.416	0.423	-1.7	91	0.00
25		Isophorone	0.838	0.775	7.5	84	0.00
26	C	2-Nitrophenol	0.193	0.189	2.1	86	0.00
27		2,4-Dimethylphenol	0.314	0.310	1.3	89	0.00
28		bis(2-Chloroethoxy)methane	0.383	0.371	3.1	85	0.00
29	C	2,4-Dichlorophenol	0.306	0.312	-2.0	90	0.00
30		1,2,4-Trichlorobenzene	0.359	0.347	3.3	88	0.00
31		Naphthalene	1.045	1.019	2.5	88	0.00
32		Benzoic acid	0.198	0.172	13.1	76	0.00
33		4-Chloroaniline	0.464	0.440	5.2	85	-0.01
34	C	Hexachlorobutadiene	0.234	0.235	-0.4	91	0.00
35		Caprolactam	0.168	0.147	12.5	74	-0.01
36	C	4-Chloro-3-methylphenol	0.398	0.386	3.0	85	0.00
37		2-Methylnaphthalene	0.803	0.793	1.2	88	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
39		1,2,4,5-Tetrachlorobenzene	0.609	0.588	3.4	88	0.00
40	P	Hexachlorocyclopentadiene	0.304	0.293	3.6	81	0.00
41	S	2,4,6-Tribromophenol	0.278	0.281	-1.1	88	0.00
42	C	2,4,6-Trichlorophenol	0.436	0.422	3.2	81	-0.01
43		2,4,5-Trichlorophenol	0.477	0.472	1.0	88	-0.01
44	S	2-Fluorobiphenyl	1.326	1.326	0.0	88	0.00

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Quant Time: Jun 05 04:00:56 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 05 03:51:43 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)
45		1,1'-Biphenyl	1.462	1.458	0.3	87	0.00
46		2-Chloronaphthalene	1.222	1.218	0.3	84	0.00
47		2-Nitroaniline	0.451	0.457	-1.3	86	0.00
48		Acenaphthylene	2.153	2.180	-1.3	88	0.00
49		Dimethylphthalate	1.741	1.626	6.6	81	0.00
50		2,6-Dinitrotoluene	0.391	0.381	2.6	85	0.00
51	C	Acenaphthene	1.269	1.265	0.3	88	0.00
52		3-Nitroaniline	0.424	0.409	3.5	84	0.00
53	P	2,4-Dinitrophenol	0.205	0.198	3.4	81	0.00
54		Dibenzofuran	1.873	1.896	-1.2	87	0.00
55	P	4-Nitrophenol	0.306	0.317	-3.6	81	0.00
56		2,4-Dinitrotoluene	0.557	0.567	-1.8	85	0.00
57		Fluorene	1.698	1.689	0.5	86	0.00
58		2,3,4,6-Tetrachlorophenol	0.435	0.435	0.0	83	0.00
59		Diethylphthalate	1.890	1.783	5.7	82	-0.01
60		4-Chlorophenyl-phenylether	0.860	0.850	1.2	86	0.00
61		4-Nitroaniline	0.466	0.492	-5.6	86	0.00
62		Azobenzene	1.796	1.890	-5.2	91	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	84	0.00
64		4,6-Dinitro-2-methylphenol	0.148	0.147	0.7	82	0.00
65	c	n-Nitrosodiphenylamine	0.597	0.612	-2.5	84	0.00
66		4-Bromophenyl-phenylether	0.221	0.228	-3.2	88	0.00
67		Hexachlorobenzene	0.238	0.244	-2.5	86	0.00
68		Atrazine	0.263	0.264	-0.4	82	0.00
69	C	Pentachlorophenol	0.128	0.133	-3.9	83	0.00
70		Phenanthrene	1.115	1.163	-4.3	88	0.00
71		Anthracene	1.114	1.163	-4.4	88	0.00
72		Carbazole	1.100	1.127	-2.5	86	0.00
73		Di-n-butylphthalate	1.409	1.417	-0.6	84	0.00
74	C	Fluoranthene	1.458	1.495	-2.5	87	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	85	0.00
76		Benzidine	0.687	0.714	-3.9	82	0.00
77		Pyrene	1.232	1.276	-3.6	86	0.00
78	S	Terphenyl-d14	0.968	1.029	-6.3	89	0.00
79		Butylbenzylphthalate	0.533	0.552	-3.6	85	0.00
80		Benzo(a)anthracene	1.230	1.279	-4.0	87	0.00
81		3,3'-Dichlorobenzidine	0.460	0.480	-4.3	86	0.00
82		Chrysene	1.114	1.159	-4.0	87	0.00
83		Bis(2-ethylhexyl)phthalate	0.769	0.815	-6.0	88	0.00
84	c	Di-n-octyl phthalate	1.289	1.367	-6.1	90	0.00
85		Indeno(1,2,3-cd)pyrene	1.399	1.473	-5.3	88	0.00
86	I	Perylene-d12	1.000	1.000	0.0	88	0.00
87		Benzo(b)fluoranthene	1.267	1.268	-0.1	85	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.213	1.238	-2.1	92	0.00
89 C	Benzo(a)pyrene	1.162	1.173	-0.9	88	0.00
90	Dibenzo(a,h)anthracene	1.208	1.220	-1.0	87	0.00
91	Benzo(g,h,i)perylene	1.150	1.160	-0.9	89	0.00

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