Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110116\
Data File : BG024591.D
Acq On : 1 Nov 2016 14:02
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 02 03:10:48 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 26 11:14:13 2016
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	137	0.00
2		1,4-Dioxane	0.499	0.476	4.6	136	0.00
3		Pyridine	1.493	1.448	3.0	138	0.00
4		n-Nitrosodimethylamine	0.773	0.693	10.3	125	0.00
5	S	2-Fluorophenol	1.220	1.159	5.0	137	0.00
6		Aniline	2.121	2.067	2.5	136	0.00
7	S	Phenol-d6	1.674	1.649	1.5	139	0.00
8		2-Chlorophenol	1.241	1.185	4.5	139	0.00
9		Benzaldehyde	1.034	0.995	3.8	136	0.00
10	C	Phenol	1.725	1.715	0.6	141	0.00
11		bis(2-Chloroethyl)ether	1.296	1.267	2.2	142	0.00
12		1,3-Dichlorobenzene	1.536	1.460	4.9	139	0.00
13	C	1,4-Dichlorobenzene	1.517	1.496	1.4	142	0.00
14		1,2-Dichlorobenzene	1.459	1.408	3.5	139	0.00
15		Benzyl Alcohol	1.557	1.496	3.9	136	0.00
16		2,2'-oxybis(1-Chloropropane	2.405	2.257	6.2	134	-0.01
17		2-Methylphenol	1.143	1.128	1.3	141	0.00
18		Hexachloroethane	0.583	0.567	2.7	145	0.00
19	P	n-Nitroso-di-n-propylamine	1.234	1.208	2.1	135	0.00
20		3+4-Methylphenols	1.535	1.559	-1.6	140	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	133	0.00
22		Acetophenone	0.514	0.504	1.9	137	0.00
23	S	Nitrobenzene-d5	0.439	0.439	0.0	141	0.00
24		Nitrobenzene	0.489	0.473	3.3	135	0.00
25		Isophorone	0.817	0.785	3.9	132	0.00
26	C	2-Nitrophenol	0.181	0.188	-3.9	145	0.00
27		2,4-Dimethylphenol	0.318	0.313	1.6	134	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.419	0.9	142	0.00
29	C	2,4-Dichlorophenol	0.333	0.328	1.5	135	0.00
30		1,2,4-Trichlorobenzene	0.395	0.395	0.0	142	0.00
31		Naphthalene	0.998	0.981	1.7	137	0.00
32		Benzoic acid	0.264	0.194	26.5#	103	0.02
33		4-Chloroaniline	0.433	0.433	0.0	133	0.00
34	C	Hexachlorobutadiene	0.309	0.314	-1.6	147	0.00
35		Caprolactam	0.122	0.115	5.7	129	0.01
36	C	4-Chloro-3-methylphenol	0.402	0.396	1.5	133	0.00
37		2-Methylnaphthalene	0.728	0.718	1.4	138	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	125	0.00
39		1,2,4,5-Tetrachlorobenzene	0.675	0.673	0.3	137	0.00
40	P	Hexachlorocyclopentadiene	0.380	0.458	-20.5	161#	0.00
41	S	2,4,6-Tribromophenol	0.299	0.311	-4.0	131	0.00
42	C	2,4,6-Trichlorophenol	0.405	0.406	-0.2	135	0.00
43		2,4,5-Trichlorophenol	0.438	0.445	-1.6	135	0.00
44	S	2-Fluorobiphenyl	1.203	1.177	2.2	132	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.388	1.381	0.5	134	0.00
46		2-Chloronaphthalene	1.057	1.052	0.5	136	0.00
47		2-Nitroaniline	0.433	0.443	-2.3	127	0.00
48		Acenaphthylene	1.621	1.596	1.5	129	0.00
49		Dimethylphthalate	1.420	1.394	1.8	127	0.00
50		2,6-Dinitrotoluene	0.303	0.326	-7.6	134	0.00
51	C	Acenaphthene	1.208	1.103	8.7	121	0.00
52		3-Nitroaniline	0.314	0.318	-1.3	129	0.00
53	P	2,4-Dinitrophenol	0.220	0.195	11.4	109	0.00
54		Dibenzofuran	1.670	1.636	2.0	129	0.00
55	P	4-Nitrophenol	0.273	0.250	8.4	117	0.00
56		2,4-Dinitrotoluene	0.440	0.467	-6.1	129	0.00
57		Fluorene	1.385	1.369	1.2	128	0.00
58		2,3,4,6-Tetrachlorophenol	0.454	0.454	0.0	126	0.00
59		Diethylphthalate	1.489	1.422	4.5	123	0.00
60		4-Chlorophenyl-phenylether	0.777	0.774	0.4	132	0.00
61		4-Nitroaniline	0.343	0.347	-1.2	129	0.00
62		Azobenzene	1.485	1.420	4.4	123	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	117	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.127	8.0	112	0.00
65	c	n-Nitrosodiphenylamine	0.547	0.539	1.5	124	0.00
66		4-Bromophenyl-phenylether	0.243	0.247	-1.6	129	0.00
67		Hexachlorobenzene	0.268	0.282	-5.2	130	0.00
68		Atrazine	0.235	0.219	6.8	115	0.00
69	C	Pentachlorophenol	0.177	0.155	12.4	104	0.00
70		Phenanthrene	1.019	0.988	3.0	123	0.00
71		Anthracene	1.028	1.011	1.7	123	0.00
72		Carbazole	0.938	0.908	3.2	123	0.00
73		Di-n-butylphthalate	1.081	1.051	2.8	122	0.00
74	C	Fluoranthene	1.289	1.258	2.4	120	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	113	0.00
76		Benzidine	0.471	0.519	-10.2	131	0.00
77		Pyrene	0.978	0.987	-0.9	118	0.00
78	S	Terphenyl-d14	0.669	0.634	5.2	114	0.00
79		Butylbenzylphthalate	0.408	0.413	-1.2	121	0.00
80		Benzo(a)anthracene	1.099	1.069	2.7	119	0.00
81		3,3'-Dichlorobenzidine	0.443	0.430	2.9	117	0.00
82		Chrysene	1.046	1.026	1.9	118	0.00
83		Bis(2-ethylhexyl)phthalate	0.583	0.568	2.6	117	-0.01
84	c	Di-n-octyl phthalate	0.968	0.949	2.0	117	-0.01
85		Indeno(1,2,3-cd)pyrene	1.444	1.395	3.4	121	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	113	-0.01
87		Benzo(b)fluoranthene	1.081	1.082	-0.1	122	0.00

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88	Benzo(k)fluoranthene	1.044	1.002	4.0	116	-0.01
89 C	Benzo(a)pyrene	1.056	1.035	2.0	118	-0.01
90	Dibenzo(a,h)anthracene	1.159	1.105	4.7	120	-0.02
91	Benzo(g,h,i)perylene	1.158	1.103	4.7	121	-0.03

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