Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG110316\
Data File : BG024672.D
Acq On : 4 Nov 2016 8:14
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 04 10:45:18 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	151#	0.00
2		1,4-Dioxane	0.499	0.491	1.6	155#	0.00
3		Pyridine	1.493	1.495	-0.1	157#	-0.01
4		n-Nitrosodimethylamine	0.773	0.796	-3.0	157#	0.00
5	S	2-Fluorophenol	1.220	1.190	2.5	155#	0.00
6		Aniline	2.121	2.171	-2.4	158#	0.00
7	S	Phenol-d6	1.674	1.731	-3.4	160#	0.00
8		2-Chlorophenol	1.241	1.241	0.0	160#	0.00
9		Benzaldehyde	1.034	1.002	3.1	151#	0.00
10	C	Phenol	1.725	1.795	-4.1	163#	0.00
11		bis(2-Chloroethyl)ether	1.296	1.304	-0.6	161#	0.00
12		1,3-Dichlorobenzene	1.536	1.497	2.5	157#	0.00
13	C	1,4-Dichlorobenzene	1.517	1.537	-1.3	160#	0.00
14		1,2-Dichlorobenzene	1.459	1.436	1.6	156#	0.00
15		Benzyl Alcohol	1.557	1.608	-3.3	161#	0.00
16		2,2'-oxybis(1-Chloropropane	2.405	2.380	1.0	155#	0.00
17		2-Methylphenol	1.143	1.176	-2.9	162#	0.00
18		Hexachloroethane	0.583	0.589	-1.0	166#	0.00
19	P	n-Nitroso-di-n-propylamine	1.234	1.270	-2.9	156#	0.00
20		3+4-Methylphenols	1.535	1.623	-5.7	161#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	151#	0.00
22		Acetophenone	0.514	0.506	1.6	157#	0.00
23	S	Nitrobenzene-d5	0.439	0.442	-0.7	162#	0.00
24		Nitrobenzene	0.489	0.486	0.6	158#	0.00
25		Isophorone	0.817	0.797	2.4	153#	0.00
26	C	2-Nitrophenol	0.181	0.193	-6.6	170#	0.00
27		2,4-Dimethylphenol	0.318	0.328	-3.1	161#	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.412	2.6	159#	0.00
29	C	2,4-Dichlorophenol	0.333	0.344	-3.3	162#	0.00
30		1,2,4-Trichlorobenzene	0.395	0.401	-1.5	164#	0.00
31		Naphthalene	0.998	0.970	2.8	155#	0.00
32		Benzoic acid	0.264	0.224	15.2	135	0.04
33		4-Chloroaniline	0.433	0.449	-3.7	158#	0.00
34	C	Hexachlorobutadiene	0.309	0.310	-0.3	166#	0.00
35		Caprolactam	0.122	0.131	-7.4	167#	0.02
36	C	4-Chloro-3-methylphenol	0.402	0.413	-2.7	158#	0.00
37		2-Methylnaphthalene	0.728	0.746	-2.5	164#	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	151#	0.00
39		1,2,4,5-Tetrachlorobenzene	0.675	0.661	2.1	163#	0.00
40	P	Hexachlorocyclopentadiene	0.380	0.344	9.5	146	0.00
41	S	2,4,6-Tribromophenol	0.299	0.314	-5.0	160#	0.00
42	C	2,4,6-Trichlorophenol	0.405	0.424	-4.7	170#	0.00
43		2,4,5-Trichlorophenol	0.438	0.434	0.9	158#	0.00
44	S	2-Fluorobiphenyl	1.203	1.122	6.7	152#	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.388	1.348	2.9	158#	0.00
46		2-Chloronaphthalene	1.057	1.037	1.9	161#	0.00
47		2-Nitroaniline	0.433	0.450	-3.9	155#	0.00
48		Acenaphthylene	1.621	1.578	2.7	154#	0.00
49		Dimethylphthalate	1.420	1.386	2.4	152#	0.00
50		2,6-Dinitrotoluene	0.303	0.326	-7.6	162#	0.00
51	C	Acenaphthene	1.208	1.086	10.1	144	0.00
52		3-Nitroaniline	0.314	0.314	0.0	153#	0.00
53	P	2,4-Dinitrophenol	0.220	0.232	-5.5	156#	0.00
54		Dibenzofuran	1.670	1.624	2.8	154#	0.00
55	P	4-Nitrophenol	0.273	0.257	5.9	145	0.01
56		2,4-Dinitrotoluene	0.440	0.476	-8.2	158#	0.00
57		Fluorene	1.385	1.371	1.0	154#	0.00
58		2,3,4,6-Tetrachlorophenol	0.454	0.454	0.0	153#	0.00
59		Diethylphthalate	1.489	1.425	4.3	148	0.00
60		4-Chlorophenyl-phenylether	0.777	0.782	-0.6	161#	0.00
61		4-Nitroaniline	0.343	0.352	-2.6	157#	0.00
62		Azobenzene	1.485	1.426	4.0	149	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	140	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.142	-2.9	149	0.00
65	c	n-Nitrosodiphenylamine	0.547	0.561	-2.6	154#	0.00
66		4-Bromophenyl-phenylether	0.243	0.256	-5.3	159#	0.00
67		Hexachlorobenzene	0.268	0.285	-6.3	156#	0.00
68		Atrazine	0.235	0.230	2.1	143	0.00
69	C	Pentachlorophenol	0.177	0.208	-17.5	166#	0.00
70		Phenanthrene	1.019	1.004	1.5	148	0.00
71		Anthracene	1.028	1.002	2.5	145	0.00
72		Carbazole	0.938	0.908	3.2	146	0.00
73		Di-n-butylphthalate	1.081	1.046	3.2	144	0.00
74	C	Fluoranthene	1.289	1.235	4.2	140	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	133	0.00
76		Benzidine	0.471	0.530	-12.5	158#	0.00
77		Pyrene	0.978	0.970	0.8	136	0.00
78	S	Terphenyl-d14	0.669	0.614	8.2	130	0.00
79		Butylbenzylphthalate	0.408	0.422	-3.4	146	0.00
80		Benzo(a)anthracene	1.099	1.071	2.5	141	0.00
81		3,3'-Dichlorobenzidine	0.443	0.447	-0.9	143	0.00
82		Chrysene	1.046	1.024	2.1	139	0.00
83		Bis(2-ethylhexyl)phthalate	0.583	0.589	-1.0	143	-0.01
84	c	Di-n-octyl phthalate	0.968	0.973	-0.5	141	-0.02
85		Indeno(1,2,3-cd)pyrene	1.444	1.427	1.2	146	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	133	-0.01
87		Benzo(b)fluoranthene	1.081	1.087	-0.6	144	0.00

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88	Benzo(k)fluoranthene	1.044	1.018	2.5	139	0.00
89 C	Benzo(a)pyrene	1.056	1.052	0.4	142	0.00
90	Dibenzo(a,h)anthracene	1.159	1.129	2.6	144	-0.01
91	Benzo(g,h,i)perylene	1.158	1.113	3.9	144	-0.01

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