Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110116\
Data File : BG024607.D
Acq On : 2 Nov 2016 00:18
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 99 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 02 03:22:36 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	157#	0.00
2		1,4-Dioxane	0.499	0.481	3.6	157#	0.00
3		Pyridine	1.493	1.489	0.3	162#	0.00
4		n-Nitrosodimethylamine	0.773	0.754	2.5	155#	0.00
5	S	2-Fluorophenol	1.220	1.156	5.2	156#	0.00
6		Aniline	2.121	2.115	0.3	160#	0.00
7	S	Phenol-d6	1.674	1.687	-0.8	162#	0.00
8		2-Chlorophenol	1.241	1.253	-1.0	168#	0.00
9		Benzaldehyde	1.034	1.011	2.2	158#	0.00
10	C	Phenol	1.725	1.758	-1.9	166#	0.00
11		bis(2-Chloroethyl)ether	1.296	1.317	-1.6	169#	0.00
12		1,3-Dichlorobenzene	1.536	1.507	1.9	165#	0.00
13	C	1,4-Dichlorobenzene	1.517	1.556	-2.6	168#	0.00
14		1,2-Dichlorobenzene	1.459	1.436	1.6	163#	0.00
15		Benzyl Alcohol	1.557	1.578	-1.3	164#	0.00
16		2,2'-oxybis(1-Chloropropane	2.405	2.295	4.6	156#	0.00
17		2-Methylphenol	1.143	1.140	0.3	163#	0.00
18		Hexachloroethane	0.583	0.601	-3.1	176#	0.00
19	P	n-Nitroso-di-n-propylamine	1.234	1.220	1.1	156#	0.00
20		3+4-Methylphenols	1.535	1.609	-4.8	166#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	150#	0.00
22		Acetophenone	0.514	0.519	-1.0	160#	0.00
23	S	Nitrobenzene-d5	0.439	0.440	-0.2	161#	0.00
24		Nitrobenzene	0.489	0.485	0.8	157#	0.00
25		Isophorone	0.817	0.798	2.3	152#	0.00
26	C	2-Nitrophenol	0.181	0.196	-8.3	172#	0.00
27		2,4-Dimethylphenol	0.318	0.324	-1.9	158#	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.434	-2.6	167#	0.00
29	C	2,4-Dichlorophenol	0.333	0.347	-4.2	162#	0.00
30		1,2,4-Trichlorobenzene	0.395	0.402	-1.8	164#	0.00
31		Naphthalene	0.998	0.982	1.6	156#	0.00
32		Benzoic acid	0.264	0.216	18.2	129	0.03
33		4-Chloroaniline	0.433	0.438	-1.2	153#	0.00
34	C	Hexachlorobutadiene	0.309	0.327	-5.8	174#	0.00
35		Caprolactam	0.122	0.109	10.7	138	0.00
36	C	4-Chloro-3-methylphenol	0.402	0.397	1.2	151#	0.00
37		2-Methylnaphthalene	0.728	0.742	-1.9	162#	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	142	0.00
39		1,2,4,5-Tetrachlorobenzene	0.675	0.697	-3.3	161#	0.00
40	P	Hexachlorocyclopentadiene	0.380	0.370	2.6	147	0.00
41	S	2,4,6-Tribromophenol	0.299	0.298	0.3	142	0.00
42	C	2,4,6-Trichlorophenol	0.405	0.414	-2.2	156#	0.00
43		2,4,5-Trichlorophenol	0.438	0.441	-0.7	151#	0.00
44	S	2-Fluorobiphenyl	1.203	1.178	2.1	150	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.388	1.384	0.3	152#	0.00
46		2-Chloronaphthalene	1.057	1.049	0.8	153#	0.00
47		2-Nitroaniline	0.433	0.434	-0.2	140	0.00
48		Acenaphthylene	1.621	1.578	2.7	144	0.00
49		Dimethylphthalate	1.420	1.370	3.5	141	0.00
50		2,6-Dinitrotoluene	0.303	0.313	-3.3	146	0.00
51	C	Acenaphthene	1.208	1.065	11.8	132	0.00
52		3-Nitroaniline	0.314	0.298	5.1	136	0.00
53	P	2,4-Dinitrophenol	0.220	0.212	3.6	134	0.00
54		Dibenzofuran	1.670	1.619	3.1	144	0.00
55	P	4-Nitrophenol	0.273	0.250	8.4	132	0.00
56		2,4-Dinitrotoluene	0.440	0.439	0.2	137	0.00
57		Fluorene	1.385	1.337	3.5	141	0.00
58		2,3,4,6-Tetrachlorophenol	0.454	0.449	1.1	142	0.00
59		Diethylphthalate	1.489	1.363	8.5	133	0.00
60		4-Chlorophenyl-phenylether	0.777	0.767	1.3	148	0.00
61		4-Nitroaniline	0.343	0.330	3.8	139	0.00
62		Azobenzene	1.485	1.353	8.9	133	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	124	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.141	-2.2	132	0.00
65	c	n-Nitrosodiphenylamine	0.547	0.572	-4.6	139	0.00
66		4-Bromophenyl-phenylether	0.243	0.256	-5.3	141	0.00
67		Hexachlorobenzene	0.268	0.285	-6.3	139	0.00
68		Atrazine	0.235	0.220	6.4	121	0.00
69	C	Pentachlorophenol	0.177	0.166	6.2	118	0.00
70		Phenanthrene	1.019	0.998	2.1	131	0.00
71		Anthracene	1.028	1.007	2.0	130	0.00
72		Carbazole	0.938	0.924	1.5	132	0.00
73		Di-n-butylphthalate	1.081	1.025	5.2	125	0.00
74	C	Fluoranthene	1.289	1.234	4.3	124	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	124	0.00
76		Benzidine	0.471	0.435	7.6	121	0.00
77		Pyrene	0.978	0.938	4.1	123	0.00
78	S	Terphenyl-d14	0.669	0.604	9.7	119	0.00
79		Butylbenzylphthalate	0.408	0.403	1.2	130	0.00
80		Benzo(a)anthracene	1.099	1.073	2.4	131	0.00
81		3,3'-Dichlorobenzidine	0.443	0.438	1.1	130	0.00
82		Chrysene	1.046	1.019	2.6	129	0.00
83		Bis(2-ethylhexyl)phthalate	0.583	0.566	2.9	128	0.00
84	c	Di-n-octyl phthalate	0.968	0.956	1.2	129	-0.01
85		Indeno(1,2,3-cd)pyrene	1.444	1.430	1.0	136	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	130	-0.01
87		Benzo(b)fluoranthene	1.081	1.022	5.5	133	0.00

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88	Benzo(k)fluoranthene	1.044	0.994	4.8	133	0.00
89 C	Benzo(a)pyrene	1.056	1.001	5.2	132	0.00
90	Dibenzo(a,h)anthracene	1.159	1.073	7.4	135	-0.01
91	Benzo(g,h,i)perylene	1.158	1.073	7.3	136	-0.03

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