Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041516\
Data File : BG021705.D
Acq On : 16 Apr 2016 6:49
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 18 01:29:11 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 13 15:48:13 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	100	0.00
2		1,4-Dioxane	0.535	0.496	7.3	91	0.00
3		Pyridine	1.604	1.461	8.9	93	0.00
4		n-Nitrosodimethylamine	0.795	0.746	6.2	98	0.00
5	S	2-Fluorophenol	1.201	1.224	-1.9	103	0.01
6		Aniline	2.373	2.243	5.5	97	0.00
7	S	Phenol-d6	1.794	1.807	-0.7	102	0.02
8		2-Chlorophenol	1.440	1.450	-0.7	102	0.00
9		Benzaldehyde	1.037	0.933	10.0	95	0.00
10	C	Phenol	1.930	1.916	0.7	100	0.00
11		bis(2-Chloroethyl)ether	1.544	1.450	6.1	96	0.00
12		1,3-Dichlorobenzene	1.608	1.588	1.2	99	0.00
13	C	1,4-Dichlorobenzene	1.637	1.630	0.4	100	0.00
14		1,2-Dichlorobenzene	1.570	1.562	0.5	101	0.00
15		Benzyl Alcohol	1.371	1.345	1.9	100	0.00
16		2,2'-oxybis(1-Chloropropane	3.308	2.933	11.3	91	0.00
17		2-Methylphenol	1.334	1.317	1.3	101	0.00
18		Hexachloroethane	0.596	0.609	-2.2	103	0.00
19	P	n-Nitroso-di-n-propylamine	1.394	1.258	9.8	95	0.00
20		3+4-Methylphenols	1.826	1.867	-2.2	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.557	0.526	5.6	98	0.00
23	S	Nitrobenzene-d5	0.389	0.383	1.5	101	0.00
24		Nitrobenzene	0.417	0.408	2.2	101	0.00
25		Isophorone	0.852	0.761	10.7	97	0.00
26	C	2-Nitrophenol	0.159	0.184	-15.7	120	0.00
27		2,4-Dimethylphenol	0.361	0.323	10.5	97	0.00
28		bis(2-Chloroethoxy)methane	0.453	0.413	8.8	96	0.00
29	C	2,4-Dichlorophenol	0.308	0.322	-4.5	108	0.00
30		1,2,4-Trichlorobenzene	0.331	0.337	-1.8	104	0.00
31		Naphthalene	1.070	1.039	2.9	100	0.00
32		Benzoic acid	0.192	0.252	-31.3#	142	0.03
33		4-Chloroaniline	0.463	0.451	2.6	104	0.00
34	C	Hexachlorobutadiene	0.214	0.219	-2.3	106	0.00
35		Caprolactam	0.141	0.135	4.3	102	0.00
36	C	4-Chloro-3-methylphenol	0.387	0.396	-2.3	109	0.00
37		2-Methylnaphthalene	0.735	0.737	-0.3	105	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
39		1,2,4,5-Tetrachlorobenzene	0.625	0.628	-0.5	111	0.00
40	P	Hexachlorocyclopentadiene	0.347	0.312	10.1	95	0.00
41	S	2,4,6-Tribromophenol	0.216	0.251	-16.2	132	0.00
42	C	2,4,6-Trichlorophenol	0.399	0.425	-6.5	116	0.00
43		2,4,5-Trichlorophenol	0.430	0.451	-4.9	114	0.00
44	S	2-Fluorobiphenyl	1.365	1.308	4.2	106	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.685	1.608	4.6	105	0.00
46		2-Chloronaphthalene	1.246	1.222	1.9	109	0.00
47		2-Nitroaniline	0.427	0.441	-3.3	116	0.00
48		Acenaphthylene	2.060	2.019	2.0	110	0.00
49		Dimethylphthalate	1.729	1.605	7.2	108	0.00
50		2,6-Dinitrotoluene	0.330	0.359	-8.8	121	0.00
51	C	Acenaphthene	1.317	1.306	0.8	110	0.00
52		3-Nitroaniline	0.366	0.382	-4.4	118	0.00
53	P	2,4-Dinitrophenol	0.106	0.140	-32.1#	152#	0.00
54		Dibenzofuran	1.798	1.789	0.5	112	0.00
55	P	4-Nitrophenol	0.313	0.327	-4.5	113	0.02
56		2,4-Dinitrotoluene	0.446	0.513	-15.0	128	0.00
57		Fluorene	1.606	1.626	-1.2	115	0.00
58		2,3,4,6-Tetrachlorophenol	0.357	0.394	-10.4	125	0.00
59		Diethylphthalate	1.805	1.724	4.5	110	0.00
60		4-Chlorophenyl-phenylether	0.791	0.815	-3.0	117	0.00
61		4-Nitroaniline	0.402	0.412	-2.5	111	0.00
62		Azobenzene	1.547	1.497	3.2	110	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
64		4,6-Dinitro-2-methylphenol	0.077	0.102	-32.5#	147	0.00
65	c	n-Nitrosodiphenylamine	0.600	0.587	2.2	113	0.00
66		4-Bromophenyl-phenylether	0.214	0.221	-3.3	119	0.00
67		Hexachlorobenzene	0.226	0.237	-4.9	125	0.00
68		Atrazine	0.237	0.230	3.0	106	0.00
69	C	Pentachlorophenol	0.129	0.117	9.3	101	0.00
70		Phenanthrene	1.102	1.073	2.6	114	0.00
71		Anthracene	1.119	1.088	2.8	113	0.00
72		Carbazole	1.044	0.959	8.1	105	0.00
73		Di-n-butylphthalate	1.329	1.242	6.5	103	0.00
74	C	Fluoranthene	1.316	1.206	8.4	103	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76		Benzidine	0.623	0.561	10.0	91	0.00
77		Pyrene	1.153	1.139	1.2	103	0.00
78	S	Terphenyl-d14	0.802	0.799	0.4	104	0.00
79		Butylbenzylphthalate	0.535	0.533	0.4	100	-0.01
80		Benzo(a)anthracene	1.151	1.122	2.5	100	0.00
81		3,3'-Dichlorobenzidine	0.911	0.902	1.0	101	0.00
82		Chrysene	1.106	1.080	2.4	101	0.00
83		Bis(2-ethylhexyl)phthalate	0.783	0.764	2.4	101	-0.02
84	c	Di-n-octyl phthalate	1.312	1.307	0.4	100	-0.03
85		Indeno(1,2,3-cd)pyrene	1.315	1.354	-3.0	105	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	103	0.00
87		Benzo(b)fluoranthene	1.160	1.125	3.0	99	0.00

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88	Benzo(k)fluoranthene	1.135	1.108	2.4	103	0.00
89 C	Benzo(a)pyrene	1.124	1.114	0.9	104	0.00
90	Dibenzo(a,h)anthracene	1.127	1.141	-1.2	105	0.00
91	Benzo(g,h,i)perylene	1.106	1.105	0.1	104	0.00

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