Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011117\
Data File : BG025471.D
Acq On : 11 Jan 2017 13:21
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 11 13:57:29 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 10 11:26:22 2017
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	105	0.00
2		1,4-Dioxane	0.469	0.506	-7.9	112	0.00
3		Pyridine	1.361	1.460	-7.3	103	0.00
4		n-Nitrosodimethylamine	0.808	0.891	-10.3	106	0.00
5	S	2-Fluorophenol	1.100	1.147	-4.3	106	0.00
6		Aniline	2.100	2.233	-6.3	106	0.00
7	S	Phenol-d6	1.550	1.662	-7.2	104	0.00
8		2-Chlorophenol	1.241	1.247	-0.5	101	0.00
9		Benzaldehyde	1.134	1.103	2.7	101	0.00
10	C	Phenol	1.718	1.826	-6.3	105	0.00
11		bis(2-Chloroethyl)ether	1.324	1.369	-3.4	105	0.00
12		1,3-Dichlorobenzene	1.502	1.556	-3.6	103	0.00
13	C	1,4-Dichlorobenzene	1.557	1.592	-2.2	103	0.00
14		1,2-Dichlorobenzene	1.486	1.518	-2.2	104	0.00
15		Benzyl Alcohol	1.479	1.602	-8.3	103	0.00
16		2,2'-oxybis(1-Chloropropane	2.451	2.644	-7.9	108	0.00
17		2-Methylphenol	1.128	1.208	-7.1	106	0.00
18		Hexachloroethane	0.598	0.622	-4.0	100	0.00
19	P	n-Nitroso-di-n-propylamine	1.309	1.358	-3.7	103	0.00
20		3+4-Methylphenols	1.561	1.693	-8.5	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.556	0.574	-3.2	103	0.00
23	S	Nitrobenzene-d5	0.453	0.479	-5.7	105	0.00
24		Nitrobenzene	0.497	0.526	-5.8	106	0.00
25		Isophorone	0.820	0.858	-4.6	103	0.00
26	C	2-Nitrophenol	0.190	0.204	-7.4	106	0.00
27		2,4-Dimethylphenol	0.312	0.336	-7.7	106	0.00
28		bis(2-Chloroethoxy)methane	0.426	0.427	-0.2	105	0.00
29	C	2,4-Dichlorophenol	0.334	0.367	-9.9	109	0.00
30		1,2,4-Trichlorobenzene	0.404	0.411	-1.7	104	0.00
31		Naphthalene	0.999	1.023	-2.4	102	0.00
32		Benzoic acid	0.251	0.250	0.4	97	0.00
33		4-Chloroaniline	0.420	0.449	-6.9	103	0.00
34	C	Hexachlorobutadiene	0.329	0.351	-6.7	108	0.00
35		Caprolactam	0.126	0.129	-2.4	103	0.00
36	C	4-Chloro-3-methylphenol	0.412	0.435	-5.6	104	0.00
37		2-Methylnaphthalene	0.740	0.796	-7.6	107	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
39		1,2,4,5-Tetrachlorobenzene	0.660	0.696	-5.5	107	0.00
40	P	Hexachlorocyclopentadiene	0.192	0.207	-7.8	101	0.00
41	S	2,4,6-Tribromophenol	0.322	0.327	-1.6	101	0.00
42	C	2,4,6-Trichlorophenol	0.416	0.417	-0.2	101	0.00
43		2,4,5-Trichlorophenol	0.435	0.435	0.0	102	0.00
44	S	2-Fluorobiphenyl	1.235	1.250	-1.2	105	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.345	1.380	-2.6	105	0.00
46		2-Chloronaphthalene	1.051	1.066	-1.4	105	0.00
47		2-Nitroaniline	0.444	0.464	-4.5	103	0.00
48		Acenaphthylene	1.629	1.618	0.7	103	0.00
49		Dimethylphthalate	1.458	1.452	0.4	100	0.00
50		2,6-Dinitrotoluene	0.319	0.321	-0.6	102	0.00
51	C	Acenaphthene	1.065	1.083	-1.7	104	0.00
52		3-Nitroaniline	0.306	0.321	-4.9	104	0.00
53	P	2,4-Dinitrophenol	0.183	0.187	-2.2	98	0.00
54		Dibenzofuran	1.684	1.698	-0.8	102	0.00
55	P	4-Nitrophenol	0.229	0.231	-0.9	98	0.00
56		2,4-Dinitrotoluene	0.464	0.491	-5.8	106	0.00
57		Fluorene	1.410	1.429	-1.3	104	0.00
58		2,3,4,6-Tetrachlorophenol	0.466	0.450	3.4	98	0.00
59		Diethylphthalate	1.529	1.537	-0.5	102	0.00
60		4-Chlorophenyl-phenylether	0.799	0.788	1.4	101	0.00
61		4-Nitroaniline	0.308	0.316	-2.6	94	0.00
62		Azobenzene	1.475	1.551	-5.2	106	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
64		4,6-Dinitro-2-methylphenol	0.139	0.154	-10.8	102	0.00
65	c	n-Nitrosodiphenylamine	0.541	0.556	-2.8	102	0.00
66		4-Bromophenyl-phenylether	0.247	0.256	-3.6	102	0.00
67		Hexachlorobenzene	0.283	0.291	-2.8	104	0.00
68		Atrazine	0.237	0.203	14.3	85	0.00
69	C	Pentachlorophenol	0.147	0.141	4.1	91	0.00
70		Phenanthrene	1.031	1.069	-3.7	104	0.00
71		Anthracene	1.047	1.080	-3.2	104	0.00
72		Carbazole	0.937	0.983	-4.9	105	0.00
73		Di-n-butylphthalate	1.152	1.191	-3.4	104	0.00
74	C	Fluoranthene	1.361	1.462	-7.4	109	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	120	0.00
76		Benzidine	0.499	0.467	6.4	106	0.00
77		Pyrene	1.045	1.040	0.5	110	0.00
78	S	Terphenyl-d14	0.754	0.730	3.2	109	0.00
79		Butylbenzylphthalate	0.420	0.429	-2.1	118	0.00
80		Benzo(a)anthracene	1.116	1.153	-3.3	116	0.00
81		3,3'-Dichlorobenzidine	0.433	0.460	-6.2	119	0.00
82		Chrysene	1.070	1.097	-2.5	117	0.00
83		Bis(2-ethylhexyl)phthalate	0.584	0.608	-4.1	119	0.00
84	c	Di-n-octyl phthalate	0.970	1.040	-7.2	121	0.00
85		Indeno(1,2,3-cd)pyrene	1.360	1.492	-9.7	123	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	125	0.00
87		Benzo(b)fluoranthene	1.091	1.117	-2.4	122	0.00

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88	Benzo(k)fluoranthene	1.054	1.090	-3.4	121	0.00
89 C	Benzo(a)pyrene	1.054	1.092	-3.6	122	0.00
90	Dibenzo(a,h)anthracene	1.117	1.158	-3.7	121	0.00
91	Benzo(g,h,i)perylene	1.110	1.151	-3.7	123	0.00

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