Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012117\
Data File : BG025568.D
Acq On : 21 Jan 2017 11:23
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 23 02:26:22 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	107	-0.02
2		1,4-Dioxane	0.469	0.543	-15.8	122	-0.03
3		Pyridine	1.361	1.500	-10.2	108	-0.02
4		n-Nitrosodimethylamine	0.808	0.898	-11.1	109	-0.03
5	S	2-Fluorophenol	1.100	1.172	-6.5	111	-0.02
6		Aniline	2.100	2.267	-8.0	110	-0.02
7	S	Phenol-d6	1.550	1.685	-8.7	108	-0.02
8		2-Chlorophenol	1.241	1.282	-3.3	105	-0.02
9		Benzaldehyde	1.134	1.093	3.6	102	-0.02
10	C	Phenol	1.718	1.833	-6.7	107	-0.02
11		bis(2-Chloroethyl)ether	1.324	1.372	-3.6	107	-0.02
12		1,3-Dichlorobenzene	1.502	1.538	-2.4	104	-0.02
13	C	1,4-Dichlorobenzene	1.557	1.619	-4.0	107	-0.02
14		1,2-Dichlorobenzene	1.486	1.507	-1.4	105	-0.02
15		Benzyl Alcohol	1.479	1.659	-12.2	108	-0.02
16		2,2'-oxybis(1-Chloropropane	2.451	2.830	-15.5	117	-0.02
17		2-Methylphenol	1.128	1.233	-9.3	110	-0.02
18		Hexachloroethane	0.598	0.626	-4.7	103	-0.02
19	P	n-Nitroso-di-n-propylamine	1.309	1.444	-10.3	112	-0.02
20		3+4-Methylphenols	1.561	1.716	-9.9	110	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	111	-0.02
22		Acetophenone	0.556	0.566	-1.8	105	-0.02
23	S	Nitrobenzene-d5	0.453	0.473	-4.4	107	-0.02
24		Nitrobenzene	0.497	0.516	-3.8	108	-0.02
25		Isophorone	0.820	0.879	-7.2	109	-0.02
26	C	2-Nitrophenol	0.190	0.187	1.6	100	-0.02
27		2,4-Dimethylphenol	0.312	0.314	-0.6	103	-0.02
28		bis(2-Chloroethoxy)methane	0.426	0.436	-2.3	111	-0.02
29	C	2,4-Dichlorophenol	0.334	0.347	-3.9	107	-0.02
30		1,2,4-Trichlorobenzene	0.404	0.401	0.7	106	-0.02
31		Naphthalene	0.999	1.027	-2.8	106	-0.02
32		Benzoic acid	0.251	0.217	13.5	87	-0.02
33		4-Chloroaniline	0.420	0.444	-5.7	105	-0.02
34	C	Hexachlorobutadiene	0.329	0.314	4.6	100	-0.02
35		Caprolactam	0.126	0.122	3.2	101	-0.03
36	C	4-Chloro-3-methylphenol	0.412	0.408	1.0	101	-0.02
37		2-Methylnaphthalene	0.740	0.777	-5.0	108	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	108	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.660	0.678	-2.7	105	-0.02
40	P	Hexachlorocyclopentadiene	0.192	0.222	-15.6	109	-0.02
41	S	2,4,6-Tribromophenol	0.322	0.307	4.7	95	-0.02
42	C	2,4,6-Trichlorophenol	0.416	0.414	0.5	101	-0.02
43		2,4,5-Trichlorophenol	0.435	0.419	3.7	98	-0.02
44	S	2-Fluorobiphenyl	1.235	1.256	-1.7	105	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.345	1.377	-2.4	105	-0.02
46		2-Chloronaphthalene	1.051	1.082	-2.9	107	-0.02
47		2-Nitroaniline	0.444	0.488	-9.9	109	-0.02
48		Acenaphthylene	1.629	1.684	-3.4	107	-0.02
49		Dimethylphthalate	1.458	1.486	-1.9	103	-0.02
50		2,6-Dinitrotoluene	0.319	0.332	-4.1	106	-0.02
51	C	Acenaphthene	1.065	1.086	-2.0	104	-0.02
52		3-Nitroaniline	0.306	0.327	-6.9	106	-0.02
53	P	2,4-Dinitrophenol	0.183	0.196	-7.1	102	-0.01
54		Dibenzofuran	1.684	1.707	-1.4	103	-0.02
55	P	4-Nitrophenol	0.229	0.222	3.1	94	-0.01
56		2,4-Dinitrotoluene	0.464	0.472	-1.7	102	-0.02
57		Fluorene	1.410	1.439	-2.1	105	-0.02
58		2,3,4,6-Tetrachlorophenol	0.466	0.426	8.6	93	-0.02
59		Diethylphthalate	1.529	1.570	-2.7	105	-0.02
60		4-Chlorophenyl-phenylether	0.799	0.799	0.0	102	-0.02
61		4-Nitroaniline	0.308	0.331	-7.5	99	-0.02
62		Azobenzene	1.475	1.604	-8.7	109	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	104	-0.02
64		4,6-Dinitro-2-methylphenol	0.139	0.145	-4.3	94	-0.02
65	c	n-Nitrosodiphenylamine	0.541	0.566	-4.6	103	-0.02
66		4-Bromophenyl-phenylether	0.247	0.258	-4.5	102	-0.02
67		Hexachlorobenzene	0.283	0.288	-1.8	101	-0.02
68		Atrazine	0.237	0.193	18.6	80	-0.02
69	C	Pentachlorophenol	0.147	0.127	13.6	81	-0.02
70		Phenanthrene	1.031	1.060	-2.8	102	-0.02
71		Anthracene	1.047	1.078	-3.0	103	-0.02
72		Carbazole	0.937	0.984	-5.0	104	-0.02
73		Di-n-butylphthalate	1.152	1.205	-4.6	104	-0.01
74	C	Fluoranthene	1.361	1.465	-7.6	107	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	120	-0.02
76		Benzidine	0.499	0.393	21.2	89	-0.01
77		Pyrene	1.045	1.021	2.3	107	-0.02
78	S	Terphenyl-d14	0.754	0.702	6.9	104	-0.02
79		Butylbenzylphthalate	0.420	0.425	-1.2	116	-0.01
80		Benzo(a)anthracene	1.116	1.129	-1.2	113	-0.02
81		3,3'-Dichlorobenzidine	0.433	0.452	-4.4	116	-0.02
82		Chrysene	1.070	1.083	-1.2	115	-0.02
83		Bis(2-ethylhexyl)phthalate	0.584	0.618	-5.8	121	-0.02
84	c	Di-n-octyl phthalate	0.970	1.061	-9.4	123	-0.02
85		Indeno(1,2,3-cd)pyrene	1.360	1.485	-9.2	122	-0.07
86	I	Perylene-d12	1.000	1.000	0.0	124	-0.04
87		Benzo(b)fluoranthene	1.091	1.114	-2.1	121	-0.03

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88	Benzo(k)fluoranthene	1.054	1.070	-1.5	118	-0.03
89 C	Benzo(a)pyrene	1.054	1.083	-2.8	121	-0.03
90	Dibenzo(a,h)anthracene	1.117	1.159	-3.8	121	-0.06
91	Benzo(g,h,i)perylene	1.110	1.144	-3.1	122	-0.07

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