Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050819\
Data File : BG040802.D
Acq On : 9 May 2019 00:06
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 09 05:02:09 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 24 05:35:33 2019
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	116	-0.02
2		1,4-Dioxane	0.516	0.494	4.3	116	0.00
3		Pyridine	1.496	1.549	-3.5	119	-0.02
4		n-Nitrosodimethylamine	0.830	0.829	0.1	122	-0.01
5	S	2-Fluorophenol	1.135	1.159	-2.1	122	-0.01
6		Aniline	2.315	2.280	1.5	117	-0.02
7	S	Phenol-d6	1.747	1.848	-5.8	127	-0.02
8		2-Chlorophenol	1.385	1.394	-0.6	119	-0.02
9		Benzaldehyde	1.083	1.154	-6.6	131	-0.01
10	C	Phenol	1.878	1.977	-5.3	126	-0.02
11		bis(2-Chloroethyl)ether	1.408	1.422	-1.0	121	-0.02
12		1,3-Dichlorobenzene	1.512	1.543	-2.1	122	-0.02
13	C	1,4-Dichlorobenzene	1.533	1.552	-1.2	121	-0.02
14		1,2-Dichlorobenzene	1.478	1.504	-1.8	122	-0.02
15		Benzyl Alcohol	1.818	1.774	2.4	116	-0.02
16		2,2'-oxybis(1-Chloropropane	2.804	2.380	15.1	102	-0.02
17		2-Methylphenol	1.329	1.311	1.4	119	-0.02
18		Hexachloroethane	0.629	0.605	3.8	116	-0.02
19	P	n-Nitroso-di-n-propylamine	1.573	1.486	5.5	115	-0.02
20		3+4-Methylphenols	1.851	1.917	-3.6	123	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	119	-0.02
22		Acetophenone	0.556	0.562	-1.1	120	-0.02
23	S	Nitrobenzene-d5	0.433	0.475	-9.7	131	-0.02
24		Nitrobenzene	0.462	0.490	-6.1	129	-0.02
25		Isophorone	0.964	0.928	3.7	116	-0.02
26	C	2-Nitrophenol	0.166	0.208	-25.3#	152#	-0.02
27		2,4-Dimethylphenol	0.311	0.315	-1.3	122	-0.02
28		bis(2-Chloroethoxy)methane	0.484	0.473	2.3	116	-0.02
29	C	2,4-Dichlorophenol	0.353	0.390	-10.5	130	-0.02
30		1,2,4-Trichlorobenzene	0.392	0.422	-7.7	128	-0.02
31		Naphthalene	1.063	1.076	-1.2	119	-0.02
32		Benzoic acid	0.213	0.209	1.9	121	-0.03
33		4-Chloroaniline	0.471	0.474	-0.6	122	-0.02
34	C	Hexachlorobutadiene	0.289	0.322	-11.4	133	-0.02
35		Caprolactam	0.139	0.139	0.0	117	-0.01
36	C	4-Chloro-3-methylphenol	0.445	0.470	-5.6	130	-0.01
37		2-Methylnaphthalene	0.794	0.818	-3.0	122	-0.02
38		1-Methylnaphthalene	0.777	0.813	-4.6	125	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	125	-0.02
40		1,2,4,5-Tetrachlorobenzene	0.745	0.808	-8.5	136	-0.02
41	P	Hexachlorocyclopentadiene	0.440	0.352	20.0	101	-0.02
42	S	2,4,6-Tribromophenol	0.275	0.312	-13.5	144	-0.02
43	C	2,4,6-Trichlorophenol	0.450	0.498	-10.7	143	-0.02
44		2,4,5-Trichlorophenol	0.490	0.552	-12.7	143	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.385	1.419	-2.5	128	-0.02
46		1,1'-Biphenyl	1.553	1.525	1.8	123	-0.02
47		2-Chloronaphthalene	1.215	1.242	-2.2	129	-0.02
48		2-Nitroaniline	0.433	0.501	-15.7	149	-0.01
49		Acenaphthylene	2.062	2.063	-0.0	124	-0.02
50		Dimethylphthalate	1.673	1.714	-2.5	128	-0.02
51		2,6-Dinitrotoluene	0.327	0.385	-17.7	149	-0.02
52	C	Acenaphthene	1.201	1.177	2.0	126	-0.02
53		3-Nitroaniline	0.335	0.375	-11.9	141	-0.02
54	P	2,4-Dinitrophenol	0.161	0.105	34.8#	86	-0.01
55		Dibenzofuran	1.967	2.042	-3.8	132	-0.02
56	P	4-Nitrophenol	0.290	0.252	13.1	112	-0.01
57		2,4-Dinitrotoluene	0.444	0.537	-20.9	155#	-0.01
58		Fluorene	1.590	1.620	-1.9	129	-0.02
59		2,3,4,6-Tetrachlorophenol	0.494	0.559	-13.2	142	-0.02
60		Diethylphthalate	1.766	1.736	1.7	124	-0.02
61		4-Chlorophenyl-phenylether	0.925	0.991	-7.1	136	-0.02
62		4-Nitroaniline	0.374	0.400	-7.0	136	-0.02
63		Azobenzene	1.818	1.692	6.9	117	-0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	130	-0.02
65		4,6-Dinitro-2-methylphenol	0.092	0.083	9.8	126	-0.01
66	c	n-Nitrosodiphenylamine	0.588	0.570	3.1	127	-0.02
67		4-Bromophenyl-phenylether	0.249	0.259	-4.0	138	-0.02
68		Hexachlorobenzene	0.258	0.271	-5.0	140	-0.02
69		Atrazine	0.233	0.217	6.9	119	-0.02
70	C	Pentachlorophenol	0.151	0.144	4.6	124	-0.01
71		Phenanthrene	1.077	1.070	0.6	129	-0.02
72		Anthracene	1.083	1.064	1.8	127	-0.02
73		Carbazole	0.987	0.941	4.7	124	-0.02
74		Di-n-butylphthalate	1.191	1.104	7.3	121	-0.03
75	C	Fluoranthene	1.342	1.351	-0.7	131	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	132	-0.03
77		Benzidine	0.548	0.433	21.0	99	-0.02
78		Pyrene	1.254	1.246	0.6	129	-0.02
79	S	Terphenyl-d14	0.903	0.892	1.2	128	-0.02
80		Butylbenzylphthalate	0.489	0.464	5.1	126	-0.03
81		Benzo(a)anthracene	1.264	1.245	1.5	129	-0.02
82		3,3'-Dichlorobenzidine	0.451	0.425	5.8	122	-0.03
83		Chrysene	1.202	1.212	-0.8	132	-0.02
84		Bis(2-ethylhexyl)phthalate	0.705	0.644	8.7	120	-0.04
85	c	Di-n-octyl phthalate	1.188	1.072	9.8	119	-0.06
86		Indeno(1,2,3-cd)pyrene	1.453	1.115	23.3	102	-0.09
87	I	Perylene-d12	1.000	1.000	0.0	129	-0.05

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.222	1.213	0.7	127	-0.05
89	Benzo(k)fluoranthene	1.141	1.190	-4.3	134	-0.04
90 C	Benzo(a)pyrene	1.157	1.145	1.0	126	-0.05
91	Dibenzo(a,h)anthracene	1.171	0.845	27.8#	92	-0.09
92	Benzo(g,h,i)perylene	1.165	0.905	22.3	99	-0.10

(#) = Out of Range SPCC's out = 0 CCC's out = 1