Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050819\
Data File : BG040789.D
Acq On : 8 May 2019 14:31
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 08 17:51:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 08 17:46:31 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	118	-0.01
2		1,4-Dioxane	0.516	0.468	9.3	111	0.00
3		Pyridine	1.496	1.498	-0.1	117	-0.02
4		n-Nitrosodimethylamine	0.830	0.807	2.8	120	0.00
5	S	2-Fluorophenol	1.135	1.121	1.2	120	-0.01
6		Aniline	2.315	2.232	3.6	116	-0.02
7	S	Phenol-d6	1.747	1.774	-1.5	123	-0.01
8		2-Chlorophenol	1.385	1.352	2.4	117	-0.02
9		Benzaldehyde	1.083	1.090	-0.6	125	-0.01
10	C	Phenol	1.878	1.905	-1.4	122	-0.01
11		bis(2-Chloroethyl)ether	1.408	1.359	3.5	116	-0.02
12		1,3-Dichlorobenzene	1.512	1.498	0.9	120	-0.02
13	C	1,4-Dichlorobenzene	1.533	1.512	1.4	119	-0.01
14		1,2-Dichlorobenzene	1.478	1.454	1.6	119	-0.02
15		Benzyl Alcohol	1.818	1.734	4.6	114	-0.02
16		2,2'-oxybis(1-Chloropropane	2.804	2.336	16.7	101	-0.02
17		2-Methylphenol	1.329	1.307	1.7	120	-0.01
18		Hexachloroethane	0.629	0.583	7.3	113	-0.02
19	P	n-Nitroso-di-n-propylamine	1.573	1.426	9.3	112	-0.02
20		3+4-Methylphenols	1.851	1.848	0.2	120	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	116	-0.01
22		Acetophenone	0.556	0.567	-2.0	118	-0.02
23	S	Nitrobenzene-d5	0.433	0.477	-10.2	129	-0.01
24		Nitrobenzene	0.462	0.490	-6.1	126	-0.01
25		Isophorone	0.964	0.933	3.2	114	-0.03
26	C	2-Nitrophenol	0.166	0.204	-22.9#	145	-0.02
27		2,4-Dimethylphenol	0.311	0.311	0.0	118	-0.02
28		bis(2-Chloroethoxy)methane	0.484	0.474	2.1	114	-0.01
29	C	2,4-Dichlorophenol	0.353	0.388	-9.9	126	-0.02
30		1,2,4-Trichlorobenzene	0.392	0.422	-7.7	125	-0.01
31		Naphthalene	1.063	1.088	-2.4	118	-0.02
32		Benzoic acid	0.213	0.233	-9.4	132	-0.03
33		4-Chloroaniline	0.471	0.484	-2.8	122	-0.02
34	C	Hexachlorobutadiene	0.289	0.318	-10.0	129	-0.02
35		Caprolactam	0.139	0.141	-1.4	117	-0.01
36	C	4-Chloro-3-methylphenol	0.445	0.472	-6.1	127	-0.01
37		2-Methylnaphthalene	0.794	0.831	-4.7	121	-0.01
38		1-Methylnaphthalene	0.777	0.803	-3.3	121	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	124	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.745	0.786	-5.5	131	-0.02
41	P	Hexachlorocyclopentadiene	0.440	0.364	17.3	104	-0.02
42	S	2,4,6-Tribromophenol	0.275	0.311	-13.1	143	-0.01
43	C	2,4,6-Trichlorophenol	0.450	0.487	-8.2	138	-0.02
44		2,4,5-Trichlorophenol	0.490	0.541	-10.4	139	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.385	1.402	-1.2	125	-0.02
46		1,1'-Biphenyl	1.553	1.536	1.1	123	-0.02
47		2-Chloronaphthalene	1.215	1.219	-0.3	125	-0.02
48		2-Nitroaniline	0.433	0.491	-13.4	145	-0.01
49		Acenaphthylene	2.062	2.008	2.6	120	-0.01
50		Dimethylphthalate	1.673	1.672	0.1	123	-0.02
51		2,6-Dinitrotoluene	0.327	0.369	-12.8	141	-0.02
52	C	Acenaphthene	1.201	1.172	2.4	124	-0.01
53		3-Nitroaniline	0.335	0.367	-9.6	137	-0.01
54	P	2,4-Dinitrophenol	0.161	0.142	11.8	114	0.00
55		Dibenzofuran	1.967	2.019	-2.6	129	-0.01
56	P	4-Nitrophenol	0.290	0.265	8.6	117	0.00
57		2,4-Dinitrotoluene	0.444	0.538	-21.2	154#	-0.01
58		Fluorene	1.590	1.599	-0.6	126	-0.01
59		2,3,4,6-Tetrachlorophenol	0.494	0.550	-11.3	138	-0.01
60		Diethylphthalate	1.766	1.709	3.2	121	-0.02
61		4-Chlorophenyl-phenylether	0.925	0.974	-5.3	132	-0.02
62		4-Nitroaniline	0.374	0.392	-4.8	131	-0.01
63		Azobenzene	1.818	1.652	9.1	113	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	124	-0.02
65		4,6-Dinitro-2-methylphenol	0.092	0.101	-9.8	146	-0.01
66	c	n-Nitrosodiphenylamine	0.588	0.580	1.4	124	-0.01
67		4-Bromophenyl-phenylether	0.249	0.265	-6.4	135	-0.02
68		Hexachlorobenzene	0.258	0.277	-7.4	137	-0.01
69		Atrazine	0.233	0.223	4.3	117	-0.02
70	C	Pentachlorophenol	0.151	0.152	-0.7	126	-0.01
71		Phenanthrene	1.077	1.083	-0.6	125	-0.01
72		Anthracene	1.083	1.087	-0.4	125	-0.02
73		Carbazole	0.987	0.951	3.6	120	-0.01
74		Di-n-butylphthalate	1.191	1.134	4.8	119	-0.03
75	C	Fluoranthene	1.342	1.363	-1.6	127	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	128	-0.03
77		Benzidine	0.548	0.433	21.0	96	-0.01
78		Pyrene	1.254	1.246	0.6	125	-0.01
79	S	Terphenyl-d14	0.903	0.893	1.1	124	-0.02
80		Butylbenzylphthalate	0.489	0.467	4.5	122	-0.02
81		Benzo(a)anthracene	1.264	1.242	1.7	125	-0.02
82		3,3'-Dichlorobenzidine	0.451	0.420	6.9	117	-0.03
83		Chrysene	1.202	1.193	0.7	126	-0.02
84		Bis(2-ethylhexyl)phthalate	0.705	0.643	8.8	116	-0.04
85	c	Di-n-octyl phthalate	1.188	1.083	8.8	116	-0.06
86		Indeno(1,2,3-cd)pyrene	1.453	1.115	23.3	99	-0.08
87	I	Perylene-d12	1.000	1.000	0.0	122	-0.04

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.222	1.231	-0.7	122	-0.04
89	Benzo(k)fluoranthene	1.141	1.214	-6.4	130	-0.04
90 C	Benzo(a)pyrene	1.157	1.167	-0.9	122	-0.05
91	Dibenzo(a,h)anthracene	1.171	0.892	23.8	92	-0.09
92	Benzo(g,h,i)perylene	1.165	0.963	17.3	100	-0.10

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