Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122822\
Data File : BG056138.D
Acq On : 28 Dec 2022 13:13
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Dec 29 03:36:42 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 28 05:20:14 2022
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	106	0.00
2		1,4-Dioxane	0.512	0.546	-6.6	107	0.00
3		Pyridine	1.558	1.695	-8.8	111	0.00
4		n-Nitrosodimethylamine	0.317	0.327	-3.2	103	-0.01
5	S	2-Fluorophenol	1.120	1.177	-5.1	110	0.00
6		Aniline	2.256	2.285	-1.3	105	0.00
7	S	Phenol-d6	1.720	1.747	-1.6	103	0.00
8		2-Chlorophenol	1.115	1.175	-5.4	109	0.00
9		Benzaldehyde	1.023	1.045	-2.2	110	0.00
10	C	Phenol	1.745	1.783	-2.2	105	0.00
11		bis(2-Chloroethyl)ether	1.182	1.225	-3.6	108	0.00
12		1,3-Dichlorobenzene	1.357	1.367	-0.7	107	0.00
13	C	1,4-Dichlorobenzene	1.362	1.407	-3.3	107	0.00
14		1,2-Dichlorobenzene	1.314	1.312	0.2	104	0.00
15		Benzyl Alcohol	1.252	1.268	-1.3	102	0.00
16		2,2'-oxybis(1-Chloropropane	0.209	0.220	-5.3	111	0.00
17		2-Methylphenol	1.280	1.307	-2.1	104	0.00
18		Hexachloroethane	0.414	0.418	-1.0	106	0.00
19	P	n-Nitroso-di-n-propylamine	1.046	1.035	1.1	103	0.00
20		3+4-Methylphenols	1.800	1.790	0.6	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
22		Acetophenone	0.566	0.586	-3.5	103	0.00
23	S	Nitrobenzene-d5	0.391	0.401	-2.6	103	0.00
24		Nitrobenzene	0.387	0.405	-4.7	106	0.00
25		Isophorone	0.823	0.840	-2.1	103	0.00
26	C	2-Nitrophenol	0.198	0.202	-2.0	102	0.00
27		2,4-Dimethylphenol	0.362	0.364	-0.6	101	0.00
28		bis(2-Chloroethoxy)methane	0.428	0.446	-4.2	106	0.00
29	C	2,4-Dichlorophenol	0.400	0.406	-1.5	99	0.00
30		1,2,4-Trichlorobenzene	0.454	0.452	0.4	101	0.00
31		Naphthalene	1.053	1.089	-3.4	105	0.00
32		Benzoic acid	0.268	0.290	-8.2	104	0.00
33		4-Chloroaniline	0.503	0.516	-2.6	103	0.00
34	C	Hexachlorobutadiene	0.326	0.326	0.0	99	0.00
35		Caprolactam	0.135	0.146	-8.1	109	0.00
36	C	4-Chloro-3-methylphenol	0.406	0.416	-2.5	103	0.00
37		2-Methylnaphthalene	0.889	0.901	-1.3	101	0.00
38		1-Methylnaphthalene	0.825	0.816	1.1	100	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
40		1,2,4,5-Tetrachlorobenzene	0.750	0.755	-0.7	101	0.00
41	P	Hexachlorocyclopentadiene	0.427	0.422	1.2	94	0.00
42	S	2,4,6-Tribromophenol	0.253	0.254	-0.4	101	0.00
43	C	2,4,6-Trichlorophenol	0.496	0.503	-1.4	99	0.00
44		2,4,5-Trichlorophenol	0.586	0.589	-0.5	101	0.00
45	S	2-Fluorobiphenyl	1.448	1.445	0.2	100	0.00
46		1,1'-Biphenyl	1.432	1.447	-1.0	101	0.00
47		2-Chloronaphthalene	1.131	1.134	-0.3	101	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.241	0.257	-6.6	107	0.00
49		Acenaphthylene	1.840	1.841	-0.1	100	0.00
50		Dimethylphthalate	1.615	1.620	-0.3	102	0.00
51		2,6-Dinitrotoluene	0.344	0.346	-0.6	104	0.00
52	C	Acenaphthene	1.197	1.217	-1.7	102	0.00
53		3-Nitroaniline	0.327	0.341	-4.3	104	0.00
54	P	2,4-Dinitrophenol	0.246	0.249	-1.2	97	0.00
55		Dibenzofuran	1.980	2.006	-1.3	101	0.00
56	P	4-Nitrophenol	0.252	0.265	-5.2	104	0.00
57		2,4-Dinitrotoluene	0.481	0.496	-3.1	102	0.00
58		Fluorene	1.593	1.568	1.6	100	0.00
59		2,3,4,6-Tetrachlorophenol	0.551	0.548	0.5	100	0.00
60		Diethylphthalate	1.514	1.532	-1.2	102	0.00
61		4-Chlorophenyl-phenylether	0.994	0.984	1.0	99	0.00
62		4-Nitroaniline	0.335	0.355	-6.0	105	0.00
63		Azobenzene	1.168	1.187	-1.6	104	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
65		4,6-Dinitro-2-methylphenol	0.135	0.138	-2.2	100	0.00
66	c	n-Nitrosodiphenylamine	0.555	0.552	0.5	101	0.00
67		4-Bromophenyl-phenylether	0.255	0.249	2.4	99	0.00
68		Hexachlorobenzene	0.238	0.235	1.3	100	0.00
69		Atrazine	0.230	0.236	-2.6	102	0.00
70	C	Pentachlorophenol	0.183	0.191	-4.4	103	0.00
71		Phenanthrene	1.036	1.033	0.3	101	0.00
72		Anthracene	1.069	1.069	0.0	101	0.00
73		Carbazole	0.899	0.918	-2.1	103	0.00
74		Di-n-butylphthalate	0.896	0.961	-7.3	108	0.00
75	C	Fluoranthene	1.352	1.372	-1.5	103	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
77		Benzidine	0.701	0.737	-5.1	105	0.00
78		Pyrene	1.410	1.397	0.9	104	0.00
79	S	Terphenyl-d14	1.117	1.095	2.0	102	0.00
80		Butylbenzylphthalate	0.404	0.413	-2.2	105	0.00
81		Benzo(a)anthracene	1.405	1.410	-0.4	104	0.00
82		3,3'-Dichlorobenzidine	0.506	0.513	-1.4	102	0.00
83		Chrysene	1.316	1.313	0.2	103	0.00
84		Bis(2-ethylhexyl)phthalate	0.583	0.600	-2.9	106	0.00
85	c	Di-n-octyl phthalate	0.984	1.016	-3.3	106	0.00
86	I	Perylene-d12	1.000	1.000	0.0	104	0.01
87		Indeno(1,2,3-cd)pyrene	1.455	1.460	-0.3	102	0.00
88		Benzo(b)fluoranthene	1.271	1.270	0.1	102	0.00
89		Benzo(k)fluoranthene	1.265	1.263	0.2	102	0.00
90	C	Benzo(a)pyrene	1.239	1.229	0.8	101	0.00
91		Dibenzo(a,h)anthracene	1.219	1.214	0.4	102	0.00
92		Benzo(g,h,i)perylene	1.183	1.182	0.1	101	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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