Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060623\
Data File : BG057636.D
Acq On : 6 Jun 2023 14:00
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jun 06 14:34:06 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG052723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 01 03:51:58 2023
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	165#	0.00
2		1,4-Dioxane	0.477	0.462	3.1	166#	0.00
3		Pyridine	1.458	1.481	-1.6	163#	0.00
4		n-Nitrosodimethylamine	0.590	0.561	4.9	153#	0.00
5	S	2-Fluorophenol	1.151	1.170	-1.7	164#	0.00
6		Aniline	2.180	2.188	-0.4	160#	0.00
7	S	Phenol-d6	1.802	1.808	-0.3	158#	0.00
8		2-Chlorophenol	1.236	1.247	-0.9	162#	0.00
9		Benzaldehyde	1.194	1.067	10.6	145	0.00
10	C	Phenol	1.705	1.709	-0.2	158#	0.00
11		bis(2-Chloroethyl)ether	1.273	1.298	-2.0	165#	0.00
12		1,3-Dichlorobenzene	1.317	1.340	-1.7	167#	0.00
13	C	1,4-Dichlorobenzene	1.340	1.348	-0.6	163#	0.00
14		1,2-Dichlorobenzene	1.307	1.323	-1.2	165#	0.00
15		Benzyl Alcohol	1.386	1.339	3.4	149	0.00
16		2,2'-oxybis(1-Chloropropane	1.581	1.671	-5.7	173#	0.00
17		2-Methylphenol	1.252	1.264	-1.0	163#	0.00
18		Hexachloroethane	0.459	0.490	-6.8	174#	0.00
19	P	n-Nitroso-di-n-propylamine	1.252	1.282	-2.4	158#	0.00
20		3+4-Methylphenols	1.758	1.761	-0.2	157#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	161#	0.00
22		Acetophenone	0.535	0.525	1.9	159#	0.00
23	S	Nitrobenzene-d5	0.428	0.422	1.4	156#	0.00
24		Nitrobenzene	0.434	0.426	1.8	156#	0.00
25		Isophorone	0.841	0.845	-0.5	159#	0.00
26	C	2-Nitrophenol	0.186	0.191	-2.7	157#	0.00
27		2,4-Dimethylphenol	0.322	0.322	0.0	157#	0.00
28		bis(2-Chloroethoxy)methane	0.436	0.456	-4.6	166#	0.00
29	C	2,4-Dichlorophenol	0.338	0.335	0.9	156#	0.00
30		1,2,4-Trichlorobenzene	0.369	0.359	2.7	159#	0.00
31		Naphthalene	1.058	1.055	0.3	160#	0.00
32		Benzoic acid	0.167	0.147	12.0	134	0.01
33		4-Chloroaniline	0.487	0.481	1.2	156#	0.00
34	C	Hexachlorobutadiene	0.255	0.253	0.8	160#	0.00
35		Caprolactam	0.140	0.121	13.6	136	0.00
36	C	4-Chloro-3-methylphenol	0.400	0.389	2.8	149	0.00
37		2-Methylnaphthalene	0.830	0.820	1.2	156#	0.00
38		1-Methylnaphthalene	0.792	0.776	2.0	156#	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	147	0.00
40		1,2,4,5-Tetrachlorobenzene	0.656	0.691	-5.3	156#	0.00
41	P	Hexachlorocyclopentadiene	0.100	0.090	10.0	129	0.00
42	S	2,4,6-Tribromophenol	0.320	0.298	6.9	135	0.00
43	C	2,4,6-Trichlorophenol	0.413	0.430	-4.1	149	0.00
44		2,4,5-Trichlorophenol	0.492	0.500	-1.6	145	0.00
45	S	2-Fluorobiphenyl	1.392	1.444	-3.7	154#	0.00
46		1,1'-Biphenyl	1.429	1.490	-4.3	153#	0.00
47		2-Chloronaphthalene	1.077	1.119	-3.9	151#	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.377	0.378	-0.3	143	0.00
49		Acenaphthylene	1.823	1.867	-2.4	150	0.00
50		Dimethylphthalate	1.648	1.615	2.0	146	0.00
51		2,6-Dinitrotoluene	0.353	0.361	-2.3	147	0.00
52	C	Acenaphthene	1.116	1.127	-1.0	148	0.00
53		3-Nitroaniline	0.368	0.369	-0.3	145	0.00
54	P	2,4-Dinitrophenol	0.173	0.159	8.1	128	0.00
55		Dibenzofuran	1.906	1.900	0.3	147	0.00
56	P	4-Nitrophenol	0.245	0.222	9.4	130	0.00
57		2,4-Dinitrotoluene	0.528	0.514	2.7	141	0.00
58		Fluorene	1.610	1.573	2.3	144	0.00
59		2,3,4,6-Tetrachlorophenol	0.461	0.434	5.9	133	0.00
60		Diethylphthalate	1.720	1.656	3.7	143	0.00
61		4-Chlorophenyl-phenylether	0.897	0.892	0.6	147	0.00
62		4-Nitroaniline	0.397	0.378	4.8	135	0.00
63		Azobenzene	1.530	1.527	0.2	146	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	135	0.00
65		4,6-Dinitro-2-methylphenol	0.122	0.126	-3.3	136	0.00
66	c	n-Nitrosodiphenylamine	0.520	0.557	-7.1	143	0.00
67		4-Bromophenyl-phenylether	0.224	0.240	-7.1	140	0.00
68		Hexachlorobenzene	0.231	0.240	-3.9	140	0.00
69		Atrazine	0.223	0.215	3.6	132	0.00
70	C	Pentachlorophenol	0.118	0.107	9.3	120	0.00
71		Phenanthrene	1.023	1.038	-1.5	136	0.00
72		Anthracene	1.050	1.066	-1.5	135	0.00
73		Carbazole	0.939	0.921	1.9	133	0.00
74		Di-n-butylphthalate	1.115	1.106	0.8	136	0.00
75	C	Fluoranthene	1.329	1.242	6.5	128	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	121	0.00
77		Benzidine	0.605	0.547	9.6	112	0.00
78		Pyrene	1.424	1.521	-6.8	126	0.00
79	S	Terphenyl-d14	1.155	1.194	-3.4	124	0.00
80		Butylbenzylphthalate	0.519	0.554	-6.7	128	0.00
81		Benzo(a)anthracene	1.388	1.416	-2.0	122	0.00
82		3,3'-Dichlorobenzidine	0.512	0.505	1.4	118	0.01
83		Chrysene	1.311	1.325	-1.1	121	0.00
84		Bis(2-ethylhexyl)phthalate	0.740	0.789	-6.6	128	0.00
85	c	Di-n-octyl phthalate	1.226	1.325	-8.1	129	0.00
86	I	Perylene-d12	1.000	1.000	0.0	123	0.00
87		Indeno(1,2,3-cd)pyrene	1.394	1.412	-1.3	123	0.02
88		Benzo(b)fluoranthene	1.231	1.220	0.9	123	0.01
89		Benzo(k)fluoranthene	1.242	1.251	-0.7	122	0.01
90	C	Benzo(a)pyrene	1.201	1.194	0.6	121	0.02
91		Dibenzo(a,h)anthracene	1.160	1.178	-1.6	123	0.02
92		Benzo(g,h,i)perylene	1.132	1.150	-1.6	124	0.03

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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