Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042922\
Data File : BG053375.D
Acq On : 29 Apr 2022 13:38
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: May 02 01:02:26 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Apr 24 01:13:30 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	117	0.00
2		1,4-Dioxane	0.557	0.609	-9.3	122	0.00
3		Pyridine	1.470	1.627	-10.7	117	0.00
4		n-Nitrosodimethylamine	0.728	0.814	-11.8	119	0.00
5	S	2-Fluorophenol	1.167	1.267	-8.6	120	0.00
6		Aniline	2.085	2.126	-2.0	110	0.00
7	S	Phenol-d6	1.799	1.907	-6.0	114	0.00
8		2-Chlorophenol	1.260	1.293	-2.6	111	0.00
9		Benzaldehyde	1.075	1.122	-4.4	128	0.00
10	C	Phenol	1.791	1.891	-5.6	117	0.00
11		bis(2-Chloroethyl)ether	1.291	1.351	-4.6	117	0.00
12		1,3-Dichlorobenzene	1.464	1.535	-4.8	116	0.00
13	C	1,4-Dichlorobenzene	1.492	1.559	-4.5	115	0.00
14		1,2-Dichlorobenzene	1.439	1.445	-0.4	113	0.00
15		Benzyl Alcohol	1.599	1.625	-1.6	109	0.00
16		2,2'-oxybis(1-Chloropropane	2.155	2.234	-3.7	111	-0.01
17		2-Methylphenol	1.212	1.253	-3.4	114	0.00
18		Hexachloroethane	0.560	0.589	-5.2	114	0.00
19	P	n-Nitroso-di-n-propylamine	1.317	1.312	0.4	108	0.00
20		3+4-Methylphenols	1.711	1.742	-1.8	110	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	109	0.00
22		Acetophenone	0.555	0.581	-4.7	110	0.00
23	S	Nitrobenzene-d5	0.493	0.508	-3.0	109	0.00
24		Nitrobenzene	0.479	0.494	-3.1	112	0.00
25		Isophorone	0.837	0.860	-2.7	107	0.00
26	C	2-Nitrophenol	0.200	0.211	-5.5	110	0.00
27		2,4-Dimethylphenol	0.287	0.300	-4.5	109	0.00
28		bis(2-Chloroethoxy)methane	0.470	0.496	-5.5	110	0.00
29	C	2,4-Dichlorophenol	0.361	0.381	-5.5	109	0.00
30		1,2,4-Trichlorobenzene	0.408	0.425	-4.2	109	0.00
31		Naphthalene	1.067	1.084	-1.6	107	0.00
32		Benzoic acid	0.155	0.178	-14.8	98	0.00
33		4-Chloroaniline	0.457	0.461	-0.9	104	0.00
34	C	Hexachlorobutadiene	0.303	0.315	-4.0	110	0.00
35		Caprolactam	0.127	0.124	2.4	100	0.00
36	C	4-Chloro-3-methylphenol	0.420	0.426	-1.4	105	0.00
37		2-Methylnaphthalene	0.790	0.801	-1.4	107	0.00
38		1-Methylnaphthalene	0.771	0.767	0.5	105	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
40		1,2,4,5-Tetrachlorobenzene	0.679	0.708	-4.3	107	0.00
41	P	Hexachlorocyclopentadiene	0.351	0.350	0.3	98	0.00
42	S	2,4,6-Tribromophenol	0.318	0.316	0.6	99	0.00
43	C	2,4,6-Trichlorophenol	0.468	0.464	0.9	99	0.00
44		2,4,5-Trichlorophenol	0.499	0.484	3.0	97	0.00
45	S	2-Fluorobiphenyl	1.450	1.477	-1.9	104	0.00
46		1,1'-Biphenyl	1.440	1.463	-1.6	104	0.00
47		2-Chloronaphthalene	1.149	1.159	-0.9	103	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.436	0.444	-1.8	100	0.00
49		Acenaphthylene	1.799	1.832	-1.8	102	0.00
50		Dimethylphthalate	1.600	1.578	1.4	100	0.00
51		2,6-Dinitrotoluene	0.334	0.333	0.3	101	0.00
52	C	Acenaphthene	1.220	1.238	-1.5	102	0.00
53		3-Nitroaniline	0.353	0.368	-4.2	103	0.00
54	P	2,4-Dinitrophenol	0.212	0.201	5.2	92	0.00
55		Dibenzofuran	1.910	1.918	-0.4	103	0.00
56	P	4-Nitrophenol	0.269	0.271	-0.7	96	0.00
57		2,4-Dinitrotoluene	0.475	0.490	-3.2	102	0.00
58		Fluorene	1.543	1.536	0.5	101	0.00
59		2,3,4,6-Tetrachlorophenol	0.485	0.488	-0.6	100	0.00
60		Diethylphthalate	1.676	1.648	1.7	100	0.00
61		4-Chlorophenyl-phenylether	0.875	0.874	0.1	101	0.00
62		4-Nitroaniline	0.372	0.387	-4.0	106	0.00
63		Azobenzene	1.690	1.709	-1.1	103	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	0.140	0.144	-2.9	101	0.00
66	c	n-Nitrosodiphenylamine	0.575	0.583	-1.4	101	0.00
67		4-Bromophenyl-phenylether	0.256	0.265	-3.5	102	0.00
68		Hexachlorobenzene	0.278	0.282	-1.4	101	0.00
69		Atrazine	0.225	0.216	4.0	98	0.00
70	C	Pentachlorophenol	0.152	0.146	3.9	90	0.00
71		Phenanthrene	1.092	1.107	-1.4	102	0.00
72		Anthracene	1.082	1.095	-1.2	102	0.00
73		Carbazole	1.052	1.078	-2.5	103	0.00
74		Di-n-butylphthalate	1.193	1.209	-1.3	102	0.00
75	C	Fluoranthene	1.400	1.428	-2.0	104	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	109	0.00
77		Benzidine	0.571	0.541	5.3	108	0.00
78		Pyrene	1.424	1.383	2.9	103	0.00
79	S	Terphenyl-d14	1.184	1.127	4.8	101	0.00
80		Butylbenzylphthalate	0.531	0.547	-3.0	108	0.00
81		Benzo(a)anthracene	1.367	1.401	-2.5	109	0.00
82		3,3'-Dichlorobenzidine	0.503	0.510	-1.4	108	0.00
83		Chrysene	1.269	1.292	-1.8	108	0.00
84		Bis(2-ethylhexyl)phthalate	0.718	0.753	-4.9	112	0.00
85	c	Di-n-octyl phthalate	1.201	1.280	-6.6	113	0.00
86	I	Perylene-d12	1.000	1.000	0.0	115	0.00
87		Indeno(1,2,3-cd)pyrene	1.486	1.529	-2.9	114	0.00
88		Benzo(b)fluoranthene	1.262	1.293	-2.5	114	-0.01
89		Benzo(k)fluoranthene	1.240	1.256	-1.3	113	0.00
90	C	Benzo(a)pyrene	1.075	1.113	-3.5	115	0.00
91		Dibenzo(a,h)anthracene	1.223	1.247	-2.0	114	0.00
92		Benzo(g,h,i)perylene	1.205	1.254	-4.1	116	-0.01

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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