Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051322\
Data File : BG053535.D
Acq On : 13 May 2022 9:42
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: May 14 03:12:17 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 11 06:00:46 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	114	0.00
2		1,4-Dioxane	0.618	0.530	14.2	99	0.00
3		Pyridine	1.510	1.502	0.5	108	0.00
4		n-Nitrosodimethylamine	0.736	0.711	3.4	108	0.00
5	S	2-Fluorophenol	1.178	1.132	3.9	109	0.00
6		Aniline	1.981	2.010	-1.5	112	0.00
7	S	Phenol-d6	1.788	1.801	-0.7	112	0.00
8		2-Chlorophenol	1.232	1.168	5.2	107	0.00
9		Benzaldehyde	1.136	1.043	8.2	111	0.00
10	C	Phenol	1.749	1.695	3.1	107	0.00
11		bis(2-Chloroethyl)ether	1.305	1.317	-0.9	111	0.00
12		1,3-Dichlorobenzene	1.453	1.374	5.4	108	0.00
13	C	1,4-Dichlorobenzene	1.467	1.393	5.0	110	0.00
14		1,2-Dichlorobenzene	1.377	1.345	2.3	111	0.00
15		Benzyl Alcohol	1.495	1.548	-3.5	115	0.00
16		2,2'-oxybis(1-Chloropropane	2.123	2.194	-3.3	117	0.00
17		2-Methylphenol	1.182	1.161	1.8	111	0.00
18		Hexachloroethane	0.557	0.527	5.4	110	0.00
19	P	n-Nitroso-di-n-propylamine	1.283	1.311	-2.2	115	0.00
20		3+4-Methylphenols	1.670	1.708	-2.3	112	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.547	0.556	-1.6	115	0.00
23	S	Nitrobenzene-d5	0.470	0.486	-3.4	114	0.00
24		Nitrobenzene	0.455	0.457	-0.4	110	0.00
25		Isophorone	0.791	0.850	-7.5	117	0.00
26	C	2-Nitrophenol	0.184	0.193	-4.9	115	0.00
27		2,4-Dimethylphenol	0.277	0.284	-2.5	113	0.00
28		bis(2-Chloroethoxy)methane	0.455	0.481	-5.7	118	0.00
29	C	2,4-Dichlorophenol	0.335	0.354	-5.7	113	0.00
30		1,2,4-Trichlorobenzene	0.380	0.399	-5.0	118	0.00
31		Naphthalene	1.017	1.035	-1.8	114	0.00
32		Benzoic acid	0.147	0.151	-2.7	108	0.00
33		4-Chloroaniline	0.426	0.453	-6.3	114	0.00
34	C	Hexachlorobutadiene	0.282	0.292	-3.5	115	0.00
35		Caprolactam	0.122	0.132	-8.2	119	0.00
36	C	4-Chloro-3-methylphenol	0.400	0.434	-8.5	116	0.00
37		2-Methylnaphthalene	0.759	0.803	-5.8	117	0.00
38		1-Methylnaphthalene	0.741	0.767	-3.5	118	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	120	0.00
40		1,2,4,5-Tetrachlorobenzene	0.630	0.639	-1.4	119	0.00
41	P	Hexachlorocyclopentadiene	0.200	0.208	-4.0	116	0.00
42	S	2,4,6-Tribromophenol	0.295	0.303	-2.7	116	0.00
43	C	2,4,6-Trichlorophenol	0.412	0.422	-2.4	115	0.00
44		2,4,5-Trichlorophenol	0.438	0.453	-3.4	116	0.00
45	S	2-Fluorobiphenyl	1.369	1.366	0.2	119	0.00
46		1,1'-Biphenyl	1.363	1.374	-0.8	119	0.00
47		2-Chloronaphthalene	1.068	1.071	-0.3	120	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.398	0.421	-5.8	120	0.00
49		Acenaphthylene	1.704	1.719	-0.9	117	0.00
50		Dimethylphthalate	1.520	1.525	-0.3	117	0.00
51		2,6-Dinitrotoluene	0.310	0.315	-1.6	117	0.00
52	C	Acenaphthene	1.016	1.032	-1.6	118	0.00
53		3-Nitroaniline	0.335	0.340	-1.5	116	0.00
54	P	2,4-Dinitrophenol	0.166	0.164	1.2	110	0.00
55		Dibenzofuran	1.824	1.823	0.1	119	0.00
56	P	4-Nitrophenol	0.239	0.234	2.1	110	0.00
57		2,4-Dinitrotoluene	0.449	0.457	-1.8	116	0.00
58		Fluorene	1.464	1.477	-0.9	118	0.00
59		2,3,4,6-Tetrachlorophenol	0.444	0.471	-6.1	121	0.00
60		Diethylphthalate	1.570	1.583	-0.8	117	0.00
61		4-Chlorophenyl-phenylether	0.817	0.858	-5.0	124	0.00
62		4-Nitroaniline	0.347	0.345	0.6	115	0.00
63		Azobenzene	1.608	1.653	-2.8	122	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	118	0.00
65		4,6-Dinitro-2-methylphenol	0.119	0.126	-5.9	116	0.00
66	c	n-Nitrosodiphenylamine	0.525	0.531	-1.1	117	0.00
67		4-Bromophenyl-phenylether	0.234	0.242	-3.4	119	0.00
68		Hexachlorobenzene	0.256	0.258	-0.8	119	0.00
69		Atrazine	0.221	0.213	3.6	113	0.00
70	C	Pentachlorophenol	0.121	0.119	1.7	111	0.00
71		Phenanthrene	1.017	1.010	0.7	117	0.00
72		Anthracene	1.000	0.996	0.4	118	0.00
73		Carbazole	0.980	0.945	3.6	113	0.00
74		Di-n-butylphthalate	1.102	1.066	3.3	113	0.00
75	C	Fluoranthene	1.317	1.252	4.9	112	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	110	0.00
77		Benzidine	0.431	0.472	-9.5	123	0.00
78		Pyrene	1.295	1.324	-2.2	111	0.00
79	S	Terphenyl-d14	1.074	1.075	-0.1	110	0.00
80		Butylbenzylphthalate	0.475	0.492	-3.6	110	0.00
81		Benzo(a)anthracene	1.276	1.287	-0.9	111	0.00
82		3,3'-Dichlorobenzidine	0.323	0.338	-4.6	111	0.00
83		Chrysene	1.191	1.193	-0.2	109	0.00
84		Bis(2-ethylhexyl)phthalate	0.655	0.680	-3.8	112	0.00
85	c	Di-n-octyl phthalate	1.108	1.167	-5.3	113	0.00
86	I	Perylene-d12	1.000	1.000	0.0	111	0.00
87		Indeno(1,2,3-cd)pyrene	1.393	1.455	-4.5	114	0.00
88		Benzo(b)fluoranthene	1.184	1.212	-2.4	110	0.00
89		Benzo(k)fluoranthene	1.163	1.177	-1.2	114	0.00
90	C	Benzo(a)pyrene	1.008	1.036	-2.8	112	0.00
91		Dibenzo(a,h)anthracene	1.148	1.181	-2.9	113	-0.01
92		Benzo(g,h,i)perylene	1.138	1.177	-3.4	113	0.00

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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