Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122722\
Data File : BF131833.D
Acq On : 27 Dec 2022 12:15
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Dec 28 00:31:25 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 26 05:13:34 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	111	0.00
2		1,4-Dioxane	0.560	0.562	-0.4	109	0.00
3		Pyridine	1.573	1.634	-3.9	109	0.00
4		n-Nitrosodimethylamine	0.740	0.718	3.0	105	0.00
5	S	2-Fluorophenol	1.207	1.215	-0.7	109	0.00
6		Aniline	1.923	1.925	-0.1	109	0.00
7	S	Phenol-d6	1.586	1.579	0.4	110	0.00
8		2-Chlorophenol	1.271	1.286	-1.2	110	0.00
9		Benzaldehyde	1.017	0.963	5.3	113	0.00
10	C	Phenol	1.888	1.890	-0.1	109	0.00
11		bis(2-Chloroethyl)ether	1.350	1.349	0.1	110	0.00
12		1,3-Dichlorobenzene	1.453	1.454	-0.1	110	0.00
13	C	1,4-Dichlorobenzene	1.464	1.468	-0.3	109	0.00
14		1,2-Dichlorobenzene	1.376	1.364	0.9	109	0.00
15		Benzyl Alcohol	1.390	1.392	-0.1	108	0.00
16		2,2'-oxybis(1-Chloropropane	1.723	1.750	-1.6	111	0.00
17		2-Methylphenol	1.162	1.169	-0.6	110	0.00
18		Hexachloroethane	0.591	0.586	0.8	108	0.00
19	P	n-Nitroso-di-n-propylamine	1.120	1.111	0.8	112	0.00
20		3+4-Methylphenols	1.517	1.518	-0.1	110	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22		Acetophenone	0.586	0.593	-1.2	112	0.00
23	S	Nitrobenzene-d5	0.501	0.498	0.6	108	0.00
24		Nitrobenzene	0.502	0.507	-1.0	109	0.00
25		Isophorone	0.818	0.832	-1.7	111	0.00
26	C	2-Nitrophenol	0.195	0.197	-1.0	109	0.00
27		2,4-Dimethylphenol	0.278	0.282	-1.4	110	0.00
28		bis(2-Chloroethoxy)methane	0.439	0.454	-3.4	114	0.00
29	C	2,4-Dichlorophenol	0.335	0.343	-2.4	112	0.00
30		1,2,4-Trichlorobenzene	0.398	0.403	-1.3	110	0.00
31		Naphthalene	1.095	1.099	-0.4	110	0.00
32		Benzoic acid	0.204	0.224	-9.8	112	0.00
33		4-Chloroaniline	0.420	0.425	-1.2	110	0.00
34	C	Hexachlorobutadiene	0.272	0.271	0.4	110	0.00
35		Caprolactam	0.099	0.103	-4.0	112	0.00
36	C	4-Chloro-3-methylphenol	0.379	0.392	-3.4	112	0.00
37		2-Methylnaphthalene	0.735	0.755	-2.7	113	0.00
38		1-Methylnaphthalene	0.714	0.736	-3.1	113	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
40		1,2,4,5-Tetrachlorobenzene	0.702	0.729	-3.8	114	0.00
41	P	Hexachlorocyclopentadiene	0.268	0.291	-8.6	113	0.00
42	S	2,4,6-Tribromophenol	0.217	0.230	-6.0	113	0.00
43	C	2,4,6-Trichlorophenol	0.450	0.468	-4.0	112	0.00
44		2,4,5-Trichlorophenol	0.486	0.505	-3.9	111	0.00
45	S	2-Fluorobiphenyl	1.460	1.483	-1.6	112	0.00
46		1,1'-Biphenyl	1.531	1.574	-2.8	113	0.00
47		2-Chloronaphthalene	1.267	1.301	-2.7	112	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.464	0.477	-2.8	110	0.00
49		Acenaphthylene	1.886	1.936	-2.7	111	0.00
50		Dimethylphthalate	1.513	1.593	-5.3	114	0.00
51		2,6-Dinitrotoluene	0.328	0.349	-6.4	113	0.00
52	C	Acenaphthene	1.138	1.182	-3.9	112	0.00
53		3-Nitroaniline	0.334	0.351	-5.1	112	0.00
54	P	2,4-Dinitrophenol	0.188	0.189	-0.5	107	0.00
55		Dibenzofuran	1.754	1.814	-3.4	112	0.00
56	P	4-Nitrophenol	0.234	0.252	-7.7	111	0.00
57		2,4-Dinitrotoluene	0.448	0.472	-5.4	112	0.00
58		Fluorene	1.432	1.483	-3.6	112	0.00
59		2,3,4,6-Tetrachlorophenol	0.399	0.429	-7.5	115	0.00
60		Diethylphthalate	1.456	1.562	-7.3	116	0.00
61		4-Chlorophenyl-phenylether	0.744	0.782	-5.1	114	0.00
62		4-Nitroaniline	0.348	0.366	-5.2	112	0.00
63		Azobenzene	1.616	1.701	-5.3	115	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	114	0.00
65		4,6-Dinitro-2-methylphenol	0.136	0.141	-3.7	112	0.00
66	c	n-Nitrosodiphenylamine	0.610	0.622	-2.0	115	0.00
67		4-Bromophenyl-phenylether	0.232	0.237	-2.2	116	0.00
68		Hexachlorobenzene	0.255	0.256	-0.4	114	0.00
69		Atrazine	0.201	0.215	-7.0	117	0.00
70	C	Pentachlorophenol	0.126	0.133	-5.6	110	0.00
71		Phenanthrene	1.100	1.089	1.0	112	0.00
72		Anthracene	1.076	1.082	-0.6	113	0.00
73		Carbazole	0.950	0.949	0.1	112	0.00
74		Di-n-butylphthalate	1.148	1.191	-3.7	116	0.00
75	C	Fluoranthene	1.281	1.272	0.7	111	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	113	0.00
77		Benzidine	0.329	0.377	-14.6	130	0.00
78		Pyrene	1.676	1.770	-5.6	113	0.00
79	S	Terphenyl-d14	1.182	1.256	-6.3	113	0.00
80		Butylbenzylphthalate	0.622	0.685	-10.1	115	0.00
81		Benzo(a)anthracene	1.448	1.555	-7.4	115	0.00
82		3,3'-Dichlorobenzidine	0.419	0.461	-10.0	118	0.00
83		Chrysene	1.361	1.482	-8.9	116	0.00
84		Bis(2-ethylhexyl)phthalate	0.757	0.857	-13.2	121	0.00
85	c	Di-n-octyl phthalate	1.019	1.187	-16.5	125	0.00
86	I	Perylene-d12	1.000	1.000	0.0	120	0.00
87		Indeno(1,2,3-cd)pyrene	1.271	1.149	9.6	106	-0.01
88		Benzo(b)fluoranthene	1.466	1.506	-2.7	118	0.00
89		Benzo(k)fluoranthene	1.417	1.510	-6.6	127	0.00
90	C	Benzo(a)pyrene	1.135	1.140	-0.4	120	0.00
91		Dibenzo(a,h)anthracene	1.041	0.962	7.6	108	-0.01
92		Benzo(g,h,i)perylene	1.046	0.950	9.2	104	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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