Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122022\
Data File : BF131717.D
Acq On : 20 Dec 2022 17:20
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Dec 21 04:40:26 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 03:12:37 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	107	0.00
2		1,4-Dioxane	0.555	0.556	-0.2	107	0.00
3		Pyridine	1.583	1.597	-0.9	107	0.00
4		n-Nitrosodimethylamine	0.716	0.711	0.7	104	0.00
5	S	2-Fluorophenol	1.188	1.161	2.3	105	0.00
6		Aniline	1.974	1.957	0.9	106	0.00
7	S	Phenol-d6	1.599	1.559	2.5	106	0.00
8		2-Chlorophenol	1.276	1.260	1.3	107	0.00
9		Benzaldehyde	1.073	0.950	11.5	113	0.00
10	C	Phenol	1.909	1.866	2.3	105	0.00
11		bis(2-Chloroethyl)ether	1.366	1.383	-1.2	108	0.00
12		1,3-Dichlorobenzene	1.442	1.439	0.2	108	0.00
13	C	1,4-Dichlorobenzene	1.474	1.446	1.9	107	0.00
14		1,2-Dichlorobenzene	1.380	1.358	1.6	108	0.00
15		Benzyl Alcohol	1.428	1.417	0.8	107	0.00
16		2,2'-oxybis(1-Chloropropane	1.742	1.787	-2.6	112	0.00
17		2-Methylphenol	1.204	1.192	1.0	107	0.00
18		Hexachloroethane	0.584	0.591	-1.2	110	0.00
19	P	n-Nitroso-di-n-propylamine	1.153	1.178	-2.2	111	0.00
20		3+4-Methylphenols	1.550	1.548	0.1	108	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
22		Acetophenone	0.584	0.571	2.2	108	0.00
23	S	Nitrobenzene-d5	0.493	0.488	1.0	108	0.00
24		Nitrobenzene	0.501	0.501	0.0	109	0.00
25		Isophorone	0.837	0.860	-2.7	114	0.00
26	C	2-Nitrophenol	0.196	0.197	-0.5	110	0.00
27		2,4-Dimethylphenol	0.279	0.281	-0.7	111	0.00
28		bis(2-Chloroethoxy)methane	0.444	0.460	-3.6	114	0.00
29	C	2,4-Dichlorophenol	0.337	0.343	-1.8	112	0.00
30		1,2,4-Trichlorobenzene	0.396	0.398	-0.5	110	0.00
31		Naphthalene	1.092	1.083	0.8	110	0.00
32		Benzoic acid	0.223	0.227	-1.8	104	0.00
33		4-Chloroaniline	0.426	0.430	-0.9	111	0.00
34	C	Hexachlorobutadiene	0.266	0.272	-2.3	112	0.00
35		Caprolactam	0.101	0.102	-1.0	111	0.00
36	C	4-Chloro-3-methylphenol	0.384	0.395	-2.9	113	0.00
37		2-Methylnaphthalene	0.745	0.751	-0.8	112	0.00
38		1-Methylnaphthalene	0.722	0.726	-0.6	112	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	118	0.00
40		1,2,4,5-Tetrachlorobenzene	0.684	0.690	-0.9	112	0.00
41	P	Hexachlorocyclopentadiene	0.310	0.335	-8.1	115	0.00
42	S	2,4,6-Tribromophenol	0.219	0.223	-1.8	117	0.00
43	C	2,4,6-Trichlorophenol	0.447	0.454	-1.6	114	0.00
44		2,4,5-Trichlorophenol	0.484	0.496	-2.5	114	0.00
45	S	2-Fluorobiphenyl	1.428	1.434	-0.4	115	0.00
46		1,1'-Biphenyl	1.504	1.532	-1.9	114	0.00
47		2-Chloronaphthalene	1.232	1.245	-1.1	113	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.444	0.447	-0.7	114	0.00
49		Acenaphthylene	1.848	1.855	-0.4	114	0.00
50		Dimethylphthalate	1.505	1.546	-2.7	119	0.00
51		2,6-Dinitrotoluene	0.324	0.332	-2.5	116	0.00
52	C	Acenaphthene	1.132	1.132	0.0	113	0.00
53		3-Nitroaniline	0.331	0.328	0.9	112	0.00
54	P	2,4-Dinitrophenol	0.194	0.199	-2.6	115	0.00
55		Dibenzofuran	1.736	1.739	-0.2	115	0.00
56	P	4-Nitrophenol	0.237	0.235	0.8	110	0.00
57		2,4-Dinitrotoluene	0.435	0.442	-1.6	118	0.00
58		Fluorene	1.403	1.392	0.8	114	0.00
59		2,3,4,6-Tetrachlorophenol	0.404	0.404	0.0	112	0.00
60		Diethylphthalate	1.450	1.522	-5.0	123	0.00
61		4-Chlorophenyl-phenylether	0.746	0.769	-3.1	120	0.00
62		4-Nitroaniline	0.333	0.329	1.2	113	0.00
63		Azobenzene	1.592	1.637	-2.8	119	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	119	0.00
65		4,6-Dinitro-2-methylphenol	0.137	0.143	-4.4	117	0.00
66	c	n-Nitrosodiphenylamine	0.620	0.623	-0.5	118	0.00
67		4-Bromophenyl-phenylether	0.236	0.245	-3.8	121	0.00
68		Hexachlorobenzene	0.260	0.261	-0.4	117	0.00
69		Atrazine	0.207	0.213	-2.9	119	0.00
70	C	Pentachlorophenol	0.138	0.151	-9.4	124	0.00
71		Phenanthrene	1.091	1.066	2.3	117	0.00
72		Anthracene	1.068	1.059	0.8	118	0.00
73		Carbazole	0.924	0.899	2.7	118	0.00
74		Di-n-butylphthalate	1.150	1.220	-6.1	131	0.00
75	C	Fluoranthene	1.215	1.197	1.5	121	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	133	0.00
77		Benzidine	0.367	0.414	-12.8	145	0.00
78		Pyrene	1.818	1.926	-5.9	123	0.00
79	S	Terphenyl-d14	1.283	1.411	-10.0	129	0.00
80		Butylbenzylphthalate	0.641	0.759	-18.4	145	0.00
81		Benzo(a)anthracene	1.440	1.528	-6.1	133	0.00
82		3,3'-Dichlorobenzidine	0.402	0.420	-4.5	138	0.00
83		Chrysene	1.360	1.441	-6.0	133	0.00
84		Bis(2-ethylhexyl)phthalate	0.765	0.969	-26.7#	161#	0.00
85	c	Di-n-octyl phthalate	1.068	1.295	-21.3#	157#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	112	0.00
87		Indeno(1,2,3-cd)pyrene	1.343	1.417	-5.5	108	0.00
88		Benzo(b)fluoranthene	1.415	1.376	2.8	115	0.00
89		Benzo(k)fluoranthene	1.381	1.455	-5.4	123	0.00
90	C	Benzo(a)pyrene	1.121	1.117	0.4	112	0.00
91		Dibenzo(a,h)anthracene	1.110	1.176	-5.9	108	0.00
92		Benzo(g,h,i)perylene	1.097	1.156	-5.4	107	0.00

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 1