Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072021\
Data File : BF124627.D
Acq On : 20 Jul 2021 12:52
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 20 13:41:02 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 20 13:36:28 2021
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	138	-0.01
2		1,4-Dioxane	0.378	0.448	-18.5	155#	-0.05
3		Pyridine	0.968	1.044	-7.9	146	-0.05
4		n-Nitrosodimethylamine	0.787	0.848	-7.8	142	-0.05
5	S	2-Fluorophenol	1.154	1.201	-4.1	144	-0.01
6		Aniline	1.480	1.453	1.8	136	-0.01
7	S	Phenol-d6	1.404	1.488	-6.0	143	-0.02
8		2-Chlorophenol	1.294	1.383	-6.9	144	-0.01
9		Benzaldehyde	0.824	0.722	12.4	132	-0.01
10	C	Phenol	1.371	1.420	-3.6	141	-0.02
11		bis(2-Chloroethyl)ether	1.063	1.130	-6.3	146	-0.01
12		1,3-Dichlorobenzene	1.416	1.489	-5.2	146	-0.01
13	C	1,4-Dichlorobenzene	1.447	1.511	-4.4	142	-0.01
14		1,2-Dichlorobenzene	1.375	1.412	-2.7	142	-0.01
15		Benzyl Alcohol	0.972	0.997	-2.6	136	-0.01
16		2,2'-oxybis(1-Chloropropane	1.832	1.897	-3.5	141	-0.01
17		2-Methylphenol	0.947	0.994	-5.0	142	-0.01
18		Hexachloroethane	0.545	0.575	-5.5	139	-0.01
19	P	n-Nitroso-di-n-propylamine	0.790	0.819	-3.7	140	-0.01
20		3+4-Methylphenols	1.247	1.321	-5.9	150#	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	142	-0.01
22		Acetophenone	0.454	0.465	-2.4	144	-0.01
23	S	Nitrobenzene-d5	0.317	0.318	-0.3	139	-0.01
24		Nitrobenzene	0.314	0.318	-1.3	137	-0.01
25		Isophorone	0.582	0.570	2.1	138	-0.01
26	C	2-Nitrophenol	0.169	0.177	-4.7	144	-0.01
27		2,4-Dimethylphenol	0.281	0.294	-4.6	147	-0.01
28		bis(2-Chloroethoxy)methane	0.339	0.336	0.9	142	-0.01
29	C	2,4-Dichlorophenol	0.290	0.301	-3.8	142	-0.01
30		1,2,4-Trichlorobenzene	0.321	0.318	0.9	141	-0.01
31		Naphthalene	1.034	1.025	0.9	141	0.00
32		Benzoic acid	0.196	0.206	-5.1	142	-0.02
33		4-Chloroaniline	0.391	0.355	9.2	125	-0.01
34	C	Hexachlorobutadiene	0.181	0.188	-3.9	144	-0.01
35		Caprolactam	0.073	0.079	-8.2	169#	-0.02
36	C	4-Chloro-3-methylphenol	0.286	0.292	-2.1	139	-0.01
37		2-Methylnaphthalene	0.698	0.689	1.3	139	-0.01
38		1-Methylnaphthalene	0.656	0.650	0.9	142	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	136	0.00
40		1,2,4,5-Tetrachlorobenzene	0.533	0.572	-7.3	143	-0.01
41	P	Hexachlorocyclopentadiene	0.333	0.338	-1.5	134	-0.01
42	S	2,4,6-Tribromophenol	0.156	0.169	-8.3	146	-0.01
43	C	2,4,6-Trichlorophenol	0.385	0.413	-7.3	141	0.00
44		2,4,5-Trichlorophenol	0.397	0.420	-5.8	146	-0.01
45	S	2-Fluorobiphenyl	1.372	1.395	-1.7	137	-0.01
46		1,1'-Biphenyl	1.519	1.536	-1.1	138	-0.01
47		2-Chloronaphthalene	1.191	1.227	-3.0	138	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.286	0.312	-9.1	135	-0.01
49		Acenaphthylene	1.858	1.899	-2.2	138	-0.01
50		Dimethylphthalate	1.391	1.420	-2.1	138	-0.01
51		2,6-Dinitrotoluene	0.274	0.309	-12.8	135	-0.01
52	C	Acenaphthene	1.188	1.215	-2.3	139	-0.01
53		3-Nitroaniline	0.306	0.304	0.7	131	-0.02
54	P	2,4-Dinitrophenol	0.143	0.156	-9.1	149	-0.01
55		Dibenzofuran	1.747	1.786	-2.2	138	-0.01
56	P	4-Nitrophenol	0.255	0.260	-2.0	130	-0.01
57		2,4-Dinitrotoluene	0.376	0.396	-5.3	136	-0.01
58		Fluorene	1.351	1.332	1.4	137	-0.01
59		2,3,4,6-Tetrachlorophenol	0.307	0.318	-3.6	139	-0.01
60		Diethylphthalate	1.455	1.424	2.1	133	-0.01
61		4-Chlorophenyl-phenylether	0.613	0.617	-0.7	135	0.00
62		4-Nitroaniline	0.309	0.322	-4.2	140	-0.02
63		Azobenzene	1.272	1.211	4.8	131	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	134	-0.01
65		4,6-Dinitro-2-methylphenol	0.114	0.124	-8.8	138	-0.02
66	c	n-Nitrosodiphenylamine	0.651	0.665	-2.2	133	-0.01
67		4-Bromophenyl-phenylether	0.207	0.209	-1.0	133	-0.01
68		Hexachlorobenzene	0.212	0.219	-3.3	131	-0.01
69		Atrazine	0.199	0.197	1.0	129	-0.01
70	C	Pentachlorophenol	0.134	0.140	-4.5	137	-0.01
71		Phenanthrene	1.091	1.072	1.7	128	-0.01
72		Anthracene	1.095	1.077	1.6	129	0.00
73		Carbazole	1.011	0.988	2.3	129	-0.01
74		Di-n-butylphthalate	1.363	1.284	5.8	122	-0.01
75	C	Fluoranthene	1.101	1.053	4.4	123	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	111	-0.01
77		Benzidine	0.748	0.764	-2.1	114	0.00
78		Pyrene	1.662	1.784	-7.3	120	-0.01
79	S	Terphenyl-d14	1.191	1.280	-7.5	123	-0.01
80		Butylbenzylphthalate	0.803	0.793	1.2	109	-0.01
81		Benzo(a)anthracene	1.326	1.328	-0.2	112	-0.01
82		3,3'-Dichlorobenzidine	0.347	0.332	4.3	102	-0.01
83		Chrysene	1.271	1.263	0.6	110	-0.01
84		Bis(2-ethylhexyl)phthalate	1.079	1.027	4.8	105	-0.01
85	c	Di-n-octyl phthalate	1.561	1.494	4.3	105	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	121	-0.01
87		Indeno(1,2,3-cd)pyrene	1.160	1.340	-15.5	133	-0.02
88		Benzo(b)fluoranthene	1.436	1.495	-4.1	123	-0.01
89		Benzo(k)fluoranthene	1.317	1.242	5.7	104	-0.01
90	C	Benzo(a)pyrene	1.185	1.219	-2.9	122	-0.01
91		Dibenzo(a,h)anthracene	0.987	1.155	-17.0	137	-0.02
92		Benzo(g,h,i)perylene	0.964	1.119	-16.1	133	-0.03

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