Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071321\
Data File : BF124558.D
Acq On : 13 Jul 2021 21:52
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 14 02:35:13 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 08 07:49:42 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	142	0.00
2		1,4-Dioxane	0.378	0.437	-15.6	155#	0.05
3		Pyridine	0.968	1.032	-6.6	148	0.04
4		n-Nitrosodimethylamine	0.787	0.842	-7.0	144	0.04
5	S	2-Fluorophenol	1.154	1.165	-1.0	143	0.00
6		Aniline	1.480	1.546	-4.5	149	0.00
7	S	Phenol-d6	1.404	1.462	-4.1	144	0.00
8		2-Chlorophenol	1.294	1.373	-6.1	146	0.00
9		Benzaldehyde	0.824	0.738	10.4	138	0.00
10	C	Phenol	1.371	1.417	-3.4	144	0.00
11		bis(2-Chloroethyl)ether	1.063	1.119	-5.3	148	0.00
12		1,3-Dichlorobenzene	1.416	1.544	-9.0	155#	0.00
13	C	1,4-Dichlorobenzene	1.447	1.541	-6.5	148	0.00
14		1,2-Dichlorobenzene	1.375	1.414	-2.8	145	0.00
15		Benzyl Alcohol	0.972	0.995	-2.4	139	0.00
16		2,2'-oxybis(1-Chloropropane	1.832	1.901	-3.8	145	0.00
17		2-Methylphenol	0.947	0.976	-3.1	143	0.00
18		Hexachloroethane	0.545	0.580	-6.4	143	0.00
19	P	n-Nitroso-di-n-propylamine	0.790	0.840	-6.3	147	0.00
20		3+4-Methylphenols	1.247	1.322	-6.0	154#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	147	0.00
22		Acetophenone	0.454	0.465	-2.4	149	0.00
23	S	Nitrobenzene-d5	0.317	0.319	-0.6	144	0.00
24		Nitrobenzene	0.314	0.319	-1.6	143	0.00
25		Isophorone	0.582	0.569	2.2	142	0.00
26	C	2-Nitrophenol	0.169	0.189	-11.8	158#	0.00
27		2,4-Dimethylphenol	0.281	0.293	-4.3	151#	0.00
28		bis(2-Chloroethoxy)methane	0.339	0.337	0.6	147	0.00
29	C	2,4-Dichlorophenol	0.290	0.294	-1.4	143	0.00
30		1,2,4-Trichlorobenzene	0.321	0.323	-0.6	147	0.00
31		Naphthalene	1.034	1.070	-3.5	152#	0.00
32		Benzoic acid	0.196	0.207	-5.6	149	0.02
33		4-Chloroaniline	0.391	0.392	-0.3	143	0.00
34	C	Hexachlorobutadiene	0.181	0.190	-5.0	150#	0.00
35		Caprolactam	0.073	0.076	-4.1	167#	0.00
36	C	4-Chloro-3-methylphenol	0.286	0.287	-0.3	141	0.00
37		2-Methylnaphthalene	0.698	0.711	-1.9	148	0.00
38		1-Methylnaphthalene	0.656	0.658	-0.3	148	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	143	0.00
40		1,2,4,5-Tetrachlorobenzene	0.533	0.581	-9.0	153#	0.00
41	P	Hexachlorocyclopentadiene	0.333	0.349	-4.8	146	0.00
42	S	2,4,6-Tribromophenol	0.156	0.162	-3.8	147	0.00
43	C	2,4,6-Trichlorophenol	0.385	0.410	-6.5	148	0.00
44		2,4,5-Trichlorophenol	0.397	0.409	-3.0	151#	0.00
45	S	2-Fluorobiphenyl	1.372	1.418	-3.4	147	0.00
46		1,1'-Biphenyl	1.519	1.593	-4.9	151#	0.00
47		2-Chloronaphthalene	1.191	1.253	-5.2	149	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.286	0.310	-8.4	141	0.00
49		Acenaphthylene	1.858	1.912	-2.9	147	0.00
50		Dimethylphthalate	1.391	1.443	-3.7	148	0.00
51		2,6-Dinitrotoluene	0.274	0.305	-11.3	141	0.00
52	C	Acenaphthene	1.188	1.264	-6.4	152#	0.00
53		3-Nitroaniline	0.306	0.334	-9.2	152#	0.00
54	P	2,4-Dinitrophenol	0.143	0.160	-11.9	161#	0.00
55		Dibenzofuran	1.747	1.807	-3.4	148	0.00
56	P	4-Nitrophenol	0.255	0.277	-8.6	147	0.00
57		2,4-Dinitrotoluene	0.376	0.424	-12.8	154#	0.00
58		Fluorene	1.351	1.368	-1.3	149	0.00
59		2,3,4,6-Tetrachlorophenol	0.307	0.326	-6.2	150	0.00
60		Diethylphthalate	1.455	1.481	-1.8	146	0.00
61		4-Chlorophenyl-phenylether	0.613	0.645	-5.2	149	0.00
62		4-Nitroaniline	0.309	0.339	-9.7	156#	0.00
63		Azobenzene	1.272	1.275	-0.2	146	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	144	0.00
65		4,6-Dinitro-2-methylphenol	0.114	0.128	-12.3	154#	0.00
66	c	n-Nitrosodiphenylamine	0.651	0.677	-4.0	145	0.00
67		4-Bromophenyl-phenylether	0.207	0.211	-1.9	144	0.00
68		Hexachlorobenzene	0.212	0.225	-6.1	145	0.00
69		Atrazine	0.199	0.198	0.5	139	0.00
70	C	Pentachlorophenol	0.134	0.148	-10.4	156#	0.00
71		Phenanthrene	1.091	1.107	-1.5	142	0.00
72		Anthracene	1.095	1.115	-1.8	144	0.00
73		Carbazole	1.011	1.031	-2.0	145	0.00
74		Di-n-butylphthalate	1.363	1.384	-1.5	142	0.00
75	C	Fluoranthene	1.101	1.118	-1.5	140	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	124	0.00
77		Benzidine	0.748	0.730	2.4	122	0.00
78		Pyrene	1.662	1.799	-8.2	135	0.00
79	S	Terphenyl-d14	1.191	1.325	-11.3	142	0.00
80		Butylbenzylphthalate	0.803	0.840	-4.6	129	0.00
81		Benzo(a)anthracene	1.326	1.389	-4.8	131	0.00
82		3,3'-Dichlorobenzidine	0.347	0.355	-2.3	122	0.00
83		Chrysene	1.271	1.344	-5.7	130	0.00
84		Bis(2-ethylhexyl)phthalate	1.079	1.113	-3.2	127	0.00
85	c	Di-n-octyl phthalate	1.561	1.578	-1.1	124	0.00
86	I	Perylene-d12	1.000	1.000	0.0	128	0.00
87		Indeno(1,2,3-cd)pyrene	1.160	1.391	-19.9	146	0.00
88		Benzo(b)fluoranthene	1.436	1.526	-6.3	133	0.00
89		Benzo(k)fluoranthene	1.317	1.328	-0.8	118	0.00
90	C	Benzo(a)pyrene	1.185	1.312	-10.7	140	0.00
91		Dibenzo(a,h)anthracene	0.987	1.167	-18.2	146	0.00
92		Benzo(g,h,i)perylene	0.964	1.126	-16.8	142	0.00

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