Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111920\
Data File : BF122376.D
Acq On : 19 Nov 2020 12:32
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 19 13:59:19 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 10 12:25:42 2020
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	101	0.00
2		1,4-Dioxane	0.597	0.578	3.2	100	0.00
3		Pyridine	1.643	1.586	3.5	98	0.00
4		n-Nitrosodimethylamine	0.774	0.666	14.0	87	0.00
5	S	2-Fluorophenol	1.303	1.274	2.2	99	-0.02
6		Aniline	2.236	2.174	2.8	99	0.00
7	S	Phenol-d6	1.704	1.658	2.7	99	-0.02
8		2-Chlorophenol	1.351	1.342	0.7	99	0.00
9		Benzaldehyde	0.961	0.885	7.9	100	0.00
10	C	Phenol	1.678	1.618	3.6	98	-0.02
11		bis(2-Chloroethyl)ether	1.334	1.255	5.9	96	0.00
12		1,3-Dichlorobenzene	1.533	1.483	3.3	98	0.00
13	C	1,4-Dichlorobenzene	1.540	1.489	3.3	99	0.00
14		1,2-Dichlorobenzene	1.438	1.402	2.5	99	0.00
15		Benzyl Alcohol	1.211	1.187	2.0	98	0.00
16		2,2'-oxybis(1-Chloropropane	1.944	1.730	11.0	91	0.00
17		2-Methylphenol	1.127	1.115	1.1	101	-0.01
18		Hexachloroethane	0.536	0.516	3.7	97	0.00
19	P	n-Nitroso-di-n-propylamine	1.018	0.923	9.3	93	0.00
20		3+4-Methylphenols	1.387	1.363	1.7	99	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.521	0.488	6.3	94	0.00
23	S	Nitrobenzene-d5	0.327	0.369	-12.8	106	0.00
24		Nitrobenzene	0.374	0.395	-5.6	101	0.00
25		Isophorone	0.728	0.683	6.2	96	0.00
26	C	2-Nitrophenol	0.126	0.190	-50.8#	132	0.00
27		2,4-Dimethylphenol	0.344	0.326	5.2	95	-0.01
28		bis(2-Chloroethoxy)methane	0.445	0.422	5.2	95	0.00
29	C	2,4-Dichlorophenol	0.304	0.318	-4.6	103	-0.01
30		1,2,4-Trichlorobenzene	0.336	0.337	-0.3	99	0.00
31		Naphthalene	1.063	1.035	2.6	98	0.00
32		Benzoic acid	0.155	0.202	-30.3#	118	-0.03
33		4-Chloroaniline	0.463	0.457	1.3	99	0.00
34	C	Hexachlorobutadiene	0.193	0.193	0.0	100	0.00
35		Caprolactam	0.096	0.104	-8.3	107	-0.01
36	C	4-Chloro-3-methylphenol	0.317	0.325	-2.5	101	-0.01
37		2-Methylnaphthalene	0.702	0.693	1.3	100	0.00
38		1-Methylnaphthalene	0.658	0.648	1.5	98	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
40		1,2,4,5-Tetrachlorobenzene	0.616	0.602	2.3	101	0.00
41	P	Hexachlorocyclopentadiene	0.360	0.369	-2.5	101	0.00
42	S	2,4,6-Tribromophenol	0.215	0.248	-15.3	114	0.00
43	C	2,4,6-Trichlorophenol	0.400	0.415	-3.7	104	0.00
44		2,4,5-Trichlorophenol	0.420	0.449	-6.9	108	-0.01
45	S	2-Fluorobiphenyl	1.280	1.208	5.6	101	0.00
46		1,1'-Biphenyl	1.592	1.528	4.0	101	0.00
47		2-Chloronaphthalene	1.264	1.210	4.3	99	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.310	0.377	-21.6	112	-0.01
49		Acenaphthylene	1.942	1.885	2.9	101	0.00
50		Dimethylphthalate	1.422	1.401	1.5	104	0.00
51		2,6-Dinitrotoluene	0.275	0.331	-20.4	115	-0.01
52	C	Acenaphthene	1.202	1.192	0.8	103	0.00
53		3-Nitroaniline	0.317	0.365	-15.1	111	-0.01
54	P	2,4-Dinitrophenol	0.084	0.134	-59.5#	173#	-0.01
55		Dibenzofuran	1.761	1.710	2.9	102	0.00
56	P	4-Nitrophenol	0.271	0.304	-12.2	114	-0.02
57		2,4-Dinitrotoluene	0.332	0.444	-33.7#	123	0.00
58		Fluorene	1.349	1.306	3.2	102	0.00
59		2,3,4,6-Tetrachlorophenol	0.349	0.377	-8.0	107	-0.01
60		Diethylphthalate	1.395	1.343	3.7	101	-0.01
61		4-Chlorophenyl-phenylether	0.637	0.633	0.6	104	-0.01
62		4-Nitroaniline	0.313	0.369	-17.9	114	-0.02
63		Azobenzene	1.463	1.342	8.3	96	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	107	-0.01
65		4,6-Dinitro-2-methylphenol	0.076	0.109	-43.4#	154#	-0.01
66	c	n-Nitrosodiphenylamine	0.681	0.644	5.4	103	0.00
67		4-Bromophenyl-phenylether	0.235	0.232	1.3	106	0.00
68		Hexachlorobenzene	0.254	0.254	0.0	108	-0.01
69		Atrazine	0.170	0.170	0.0	107	-0.01
70	C	Pentachlorophenol	0.146	0.161	-10.3	116	-0.01
71		Phenanthrene	1.135	1.071	5.6	103	-0.01
72		Anthracene	1.170	1.110	5.1	103	-0.01
73		Carbazole	1.064	1.008	5.3	103	-0.01
74		Di-n-butylphthalate	1.134	1.084	4.4	103	-0.01
75	C	Fluoranthene	1.185	1.144	3.5	104	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	98	-0.02
77		Benzidine	0.555	0.521	6.1	92	-0.02
78		Pyrene	1.405	1.436	-2.2	103	-0.02
79	S	Terphenyl-d14	0.944	0.953	-1.0	105	-0.01
80		Butylbenzylphthalate	0.505	0.567	-12.3	105	-0.02
81		Benzo(a)anthracene	1.244	1.211	2.7	98	-0.02
82		3,3'-Dichlorobenzidine	0.464	0.474	-2.2	99	-0.02
83		Chrysene	1.253	1.222	2.5	99	-0.02
84		Bis(2-ethylhexyl)phthalate	0.677	0.715	-5.6	105	-0.02
85	c	Di-n-octyl phthalate	1.148	1.202	-4.7	98	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	90	-0.04
87		Indeno(1,2,3-cd)pyrene	1.200	1.164	3.0	92	-0.06
88		Benzo(b)fluoranthene	1.295	1.336	-3.2	92	-0.02
89		Benzo(k)fluoranthene	1.263	1.223	3.2	90	-0.03
90	C	Benzo(a)pyrene	1.129	1.121	0.7	90	-0.04
91		Dibenzo(a,h)anthracene	1.005	0.963	4.2	91	-0.05
92		Benzo(g,h,i)perylene	0.978	0.958	2.0	95	-0.06

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