Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF080320\
Data File : BF121172.D
Acq On : 3 Aug 2020 10:18
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 03 11:21:38 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 29 14:55:43 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	98	0.00
2		1,4-Dioxane	0.561	0.539	3.9	97	-0.01
3		Pyridine	1.494	1.449	3.0	99	-0.02
4		n-Nitrosodimethylamine	0.639	0.590	7.7	93	-0.02
5	S	2-Fluorophenol	1.220	1.168	4.3	97	0.00
6		Aniline	2.057	1.821	11.5	90	0.00
7	S	Phenol-d6	1.576	1.502	4.7	97	0.00
8		2-Chlorophenol	1.344	1.307	2.8	98	0.00
9		Benzaldehyde	0.811	0.793	2.2	104	0.00
10	C	Phenol	1.734	1.582	8.8	92	0.00
11		bis(2-Chloroethyl)ether	1.329	1.301	2.1	99	-0.01
12		1,3-Dichlorobenzene	1.457	1.422	2.4	99	0.00
13	C	1,4-Dichlorobenzene	1.449	1.427	1.5	100	0.00
14		1,2-Dichlorobenzene	1.371	1.312	4.3	96	0.00
15		Benzyl Alcohol	1.114	1.012	9.2	92	0.00
16		2,2'-oxybis(1-Chloropropane	1.915	1.758	8.2	93	-0.01
17		2-Methylphenol	1.191	1.144	3.9	99	-0.01
18		Hexachloroethane	0.525	0.519	1.1	99	0.00
19	P	n-Nitroso-di-n-propylamine	0.896	0.809	9.7	95	-0.01
20		3+4-Methylphenols	1.515	1.310	13.5	96	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
22		Acetophenone	0.449	0.440	2.0	98	0.00
23	S	Nitrobenzene-d5	0.346	0.352	-1.7	99	-0.01
24		Nitrobenzene	0.373	0.376	-0.8	99	-0.01
25		Isophorone	0.690	0.679	1.6	98	-0.01
26	C	2-Nitrophenol	0.177	0.199	-12.4	105	0.00
27		2,4-Dimethylphenol	0.287	0.287	0.0	99	0.00
28		bis(2-Chloroethoxy)methane	0.421	0.421	0.0	100	0.00
29	C	2,4-Dichlorophenol	0.294	0.300	-2.0	99	0.00
30		1,2,4-Trichlorobenzene	0.326	0.331	-1.5	99	0.00
31		Naphthalene	1.026	1.018	0.8	98	0.00
32		Benzoic acid	0.216	0.190	12.0	77	0.00
33		4-Chloroaniline	0.458	0.451	1.5	99	-0.01
34	C	Hexachlorobutadiene	0.192	0.194	-1.0	99	-0.01
35		Caprolactam	0.094	0.095	-1.1	99	0.00
36	C	4-Chloro-3-methylphenol	0.301	0.305	-1.3	99	0.00
37		2-Methylnaphthalene	0.666	0.672	-0.9	99	0.00
38		1-Methylnaphthalene	0.631	0.631	0.0	99	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
40		1,2,4,5-Tetrachlorobenzene	0.583	0.605	-3.8	101	0.00
41	P	Hexachlorocyclopentadiene	0.322	0.306	5.0	89	0.00
42	S	2,4,6-Tribromophenol	0.219	0.224	-2.3	98	0.00
43	C	2,4,6-Trichlorophenol	0.410	0.402	2.0	95	0.00
44		2,4,5-Trichlorophenol	0.465	0.464	0.2	97	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.340	1.189	11.3	98	0.00
46		1,1'-Biphenyl	1.526	1.508	1.2	97	0.00
47		2-Chloronaphthalene	1.244	1.244	0.0	98	0.00
48		2-Nitroaniline	0.376	0.393	-4.5	101	0.00
49		Acenaphthylene	1.932	1.940	-0.4	99	0.00
50		Dimethylphthalate	1.443	1.456	-0.9	101	0.00
51		2,6-Dinitrotoluene	0.310	0.327	-5.5	101	0.00
52	C	Acenaphthene	1.229	1.131	8.0	90	0.00
53		3-Nitroaniline	0.389	0.390	-0.3	98	0.00
54	P	2,4-Dinitrophenol	0.149	0.160	-7.4	105	0.00
55		Dibenzofuran	1.777	1.701	4.3	95	0.00
56	P	4-Nitrophenol	0.295	0.253	14.2	82	0.00
57		2,4-Dinitrotoluene	0.407	0.423	-3.9	98	0.00
58		Fluorene	1.325	1.292	2.5	100	0.00
59		2,3,4,6-Tetrachlorophenol	0.354	0.334	5.6	92	0.00
60		Diethylphthalate	1.459	1.422	2.5	97	0.00
61		4-Chlorophenyl-phenylether	0.707	0.652	7.8	99	0.00
62		4-Nitroaniline	0.379	0.393	-3.7	102	0.00
63		Azobenzene	1.389	1.336	3.8	95	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
65		4,6-Dinitro-2-methylphenol	0.101	0.117	-15.8	105	0.00
66	c	n-Nitrosodiphenylamine	0.629	0.641	-1.9	96	0.00
67		4-Bromophenyl-phenylether	0.223	0.235	-5.4	100	0.00
68		Hexachlorobenzene	0.241	0.250	-3.7	99	0.00
69		Atrazine	0.156	0.163	-4.5	101	0.00
70	C	Pentachlorophenol	0.138	0.141	-2.2	90	0.00
71		Phenanthrene	1.029	1.011	1.7	95	0.00
72		Anthracene	1.033	1.036	-0.3	95	0.00
73		Carbazole	1.025	1.026	-0.1	95	0.00
74		Di-n-butylphthalate	1.183	1.135	4.1	92	0.00
75	C	Fluoranthene	1.140	1.110	2.6	93	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	83	0.00
77		Benzidine	0.526	0.555	-5.5	82	0.00
78		Pyrene	1.414	1.553	-9.8	92	0.00
79	S	Terphenyl-d14	0.920	0.928	-0.9	91	0.00
80		Butylbenzylphthalate	0.630	0.661	-4.9	86	0.00
81		Benzo(a)anthracene	1.211	1.285	-6.1	92	0.00
82		3,3'-Dichlorobenzidine	0.457	0.467	-2.2	86	0.00
83		Chrysene	1.273	1.278	-0.4	84	0.00
84		Bis(2-ethylhexyl)phthalate	0.777	0.828	-6.6	90	0.00
85	c	Di-n-octyl phthalate	1.546	1.390	10.1	74	0.00
86	I	Perylene-d12	1.000	1.000	0.0	66	0.01
87		Indeno(1,2,3-cd)pyrene	1.391	1.186	14.7	56	0.02

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88	Benzo(b)fluoranthene	1.209	1.293	-6.9	72	0.00
89	Benzo(k)fluoranthene	1.103	1.264	-14.6	73	0.00
90 C	Benzo(a)pyrene	1.103	1.163	-5.4	69	0.01
91	Dibenzo(a,h)anthracene	1.136	0.973	14.3	56	0.02
92	Benzo(g,h,i)perylene	1.120	0.952	15.0	56	0.02

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