Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010820\
Data File : BG043979.D
Acq On : 8 Jan 2020 23:57
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 09 06:15:05 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 31 13:32:23 2019
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	-0.01
2		1,4-Dioxane	0.475	0.469	1.3	96	0.00
3		Pyridine	1.403	1.406	-0.2	96	0.00
4		n-Nitrosodimethylamine	0.625	0.563	9.9	89	0.00
5	S	2-Fluorophenol	1.089	1.117	-2.6	97	0.00
6		Aniline	2.000	1.901	4.9	92	-0.01
7	S	Phenol-d6	1.523	1.511	0.8	95	0.00
8		2-Chlorophenol	1.279	1.260	1.5	96	-0.01
9		Benzaldehyde	0.974	0.815	16.3	86	0.00
10	C	Phenol	1.569	1.528	2.6	94	0.00
11		bis(2-Chloroethyl)ether	1.354	1.287	4.9	92	-0.01
12		1,3-Dichlorobenzene	1.496	1.473	1.5	96	-0.01
13	C	1,4-Dichlorobenzene	1.476	1.486	-0.7	97	-0.01
14		1,2-Dichlorobenzene	1.417	1.392	1.8	96	-0.01
15		Benzyl Alcohol	1.202	1.123	6.6	91	-0.01
16		2,2'-oxybis(1-Chloropropane	2.095	1.879	10.3	89	-0.02
17		2-Methylphenol	1.135	1.087	4.2	94	0.00
18		Hexachloroethane	0.575	0.560	2.6	95	-0.01
19	P	n-Nitroso-di-n-propylamine	1.134	1.024	9.7	88	0.00
20		3+4-Methylphenols	1.584	1.507	4.9	92	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	95	-0.01
22		Acetophenone	0.497	0.466	6.2	90	0.00
23	S	Nitrobenzene-d5	0.374	0.342	8.6	90	-0.01
24		Nitrobenzene	0.370	0.342	7.6	89	-0.01
25		Isophorone	0.701	0.649	7.4	89	0.00
26	C	2-Nitrophenol	0.175	0.180	-2.9	101	-0.01
27		2,4-Dimethylphenol	0.264	0.247	6.4	92	-0.01
28		bis(2-Chloroethoxy)methane	0.468	0.432	7.7	88	-0.01
29	C	2,4-Dichlorophenol	0.315	0.303	3.8	93	0.00
30		1,2,4-Trichlorobenzene	0.353	0.338	4.2	93	-0.01
31		Naphthalene	0.968	0.940	2.9	91	0.00
32		Benzoic acid	0.197	0.171	13.2	94	0.00
33		4-Chloroaniline	0.462	0.435	5.8	89	0.00
34	C	Hexachlorobutadiene	0.215	0.205	4.7	93	-0.01
35		Caprolactam	0.117	0.111	5.1	92	0.00
36	C	4-Chloro-3-methylphenol	0.335	0.320	4.5	92	0.00
37		2-Methylnaphthalene	0.729	0.692	5.1	91	-0.01
38		1-Methylnaphthalene	0.691	0.660	4.5	92	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	96	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.586	0.553	5.6	91	-0.01
41	P	Hexachlorocyclopentadiene	0.304	0.281	7.6	89	-0.01
42	S	2,4,6-Tribromophenol	0.209	0.212	-1.4	95	-0.01
43	C	2,4,6-Trichlorophenol	0.403	0.389	3.5	92	-0.01
44		2,4,5-Trichlorophenol	0.416	0.402	3.4	93	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.104	1.093	1.0	90	-0.01
46		1,1'-Biphenyl	1.434	1.381	3.7	90	0.00
47		2-Chloronaphthalene	1.159	1.101	5.0	90	0.00
48		2-Nitroaniline	0.373	0.348	6.7	90	0.00
49		Acenaphthylene	1.665	1.625	2.4	91	0.00
50		Dimethylphthalate	1.420	1.364	3.9	89	0.00
51		2,6-Dinitrotoluene	0.321	0.303	5.6	91	-0.01
52	C	Acenaphthene	1.168	1.112	4.8	91	-0.01
53		3-Nitroaniline	0.364	0.356	2.2	93	0.00
54	P	2,4-Dinitrophenol	0.145	0.132	9.0	88	0.00
55		Dibenzofuran	1.608	1.563	2.8	90	0.00
56	P	4-Nitrophenol	0.279	0.282	-1.1	94	0.00
57		2,4-Dinitrotoluene	0.437	0.433	0.9	94	0.00
58		Fluorene	1.274	1.237	2.9	91	-0.01
59		2,3,4,6-Tetrachlorophenol	0.358	0.354	1.1	93	0.00
60		Diethylphthalate	1.420	1.386	2.4	91	-0.01
61		4-Chlorophenyl-phenylether	0.692	0.661	4.5	92	-0.01
62		4-Nitroaniline	0.360	0.359	0.3	93	0.00
63		Azobenzene	1.317	1.248	5.2	88	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
65		4,6-Dinitro-2-methylphenol	0.097	0.097	0.0	101	0.00
66	c	n-Nitrosodiphenylamine	0.589	0.543	7.8	92	-0.01
67		4-Bromophenyl-phenylether	0.225	0.202	10.2	92	-0.01
68		Hexachlorobenzene	0.242	0.224	7.4	94	-0.01
69		Atrazine	0.191	0.178	6.8	88	0.00
70	C	Pentachlorophenol	0.134	0.126	6.0	91	0.00
71		Phenanthrene	0.959	0.921	4.0	92	-0.01
72		Anthracene	0.948	0.918	3.2	93	0.00
73		Carbazole	0.876	0.872	0.5	95	0.00
74		Di-n-butylphthalate	1.118	1.073	4.0	94	-0.01
75	C	Fluoranthene	1.097	1.049	4.4	94	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	99	0.00
77		Benzidine	0.503	0.470	6.6	83	0.00
78		Pyrene	1.305	1.229	5.8	95	-0.01
79	S	Terphenyl-d14	0.828	0.807	2.5	95	-0.01
80		Butylbenzylphthalate	0.622	0.611	1.8	97	-0.01
81		Benzo(a)anthracene	1.240	1.198	3.4	94	-0.01
82		3,3'-Dichlorobenzidine	0.490	0.468	4.5	93	-0.01
83		Chrysene	1.178	1.150	2.4	95	-0.01
84		Bis(2-ethylhexyl)phthalate	0.858	0.823	4.1	94	-0.01
85	c	Di-n-octyl phthalate	1.442	1.431	0.8	96	-0.02
86		Indeno(1,2,3-cd)pyrene	1.601	1.463	8.6	92	-0.03
87	I	Perylene-d12	1.000	1.000	0.0	99	-0.02

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88	Benzo(b)fluoranthene	1.182	1.101	6.9	93	-0.02
89	Benzo(k)fluoranthene	1.124	1.097	2.4	94	-0.02
90 C	Benzo(a)pyrene	1.116	1.059	5.1	93	-0.02
91	Dibenzo(a,h)anthracene	1.121	1.043	7.0	93	-0.03
92	Benzo(g,h,i)perylene	1.113	1.021	8.3	91	-0.03

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