Evaluate Continuing Calibration Report

Data Path : U:\HPCHEM1\BNA G\DATA\BG011018\
Data File : BG031978.D
Acq On : 10 Jan 2018 15:17
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 11 00:42:02 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 05 15:31:34 2018
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	102	0.00
2		1,4-Dioxane	0.407	0.360	11.5	94	0.00
3		Pyridine	1.088	1.066	2.0	95	0.00
4		n-Nitrosodimethylamine	0.439	0.443	-0.9	101	0.00
5	S	2-Fluorophenol	1.098	1.049	4.5	98	0.00
6		Aniline	1.834	1.735	5.4	94	0.00
7	S	Phenol-d6	1.426	1.429	-0.2	99	0.00
8		2-Chlorophenol	1.274	1.251	1.8	100	0.00
9		Benzaldehyde	0.858	0.714	16.8	83	0.00
10	C	Phenol	1.439	1.421	1.3	100	0.00
11		bis(2-Chloroethyl)ether	1.195	1.095	8.4	93	0.00
12		1,3-Dichlorobenzene	1.582	1.535	3.0	99	0.00
13	C	1,4-Dichlorobenzene	1.616	1.587	1.8	102	0.00
14		1,2-Dichlorobenzene	1.531	1.509	1.4	101	0.00
15		Benzyl Alcohol	1.070	1.113	-4.0	102	0.00
16		2,2'-oxybis(1-Chloropropane	0.973	0.851	12.5	90	-0.01
17		2-Methylphenol	1.030	1.026	0.4	99	0.00
18		Hexachloroethane	0.489	0.491	-0.4	106	0.00
19	P	n-Nitroso-di-n-propylamine	0.974	0.936	3.9	98	0.00
20		3+4-Methylphenols	1.464	1.456	0.5	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
22		Acetophenone	0.465	0.445	4.3	98	0.00
23	S	Nitrobenzene-d5	0.265	0.298	-12.5	111	0.00
24		Nitrobenzene	0.273	0.287	-5.1	103	-0.01
25		Isophorone	0.570	0.540	5.3	95	-0.01
26	C	2-Nitrophenol	0.143	0.181	-26.6#	127	0.00
27		2,4-Dimethylphenol	0.301	0.283	6.0	98	-0.01
28		bis(2-Chloroethoxy)methane	0.383	0.353	7.8	94	-0.01
29	C	2,4-Dichlorophenol	0.354	0.361	-2.0	102	0.00
30		1,2,4-Trichlorobenzene	0.432	0.415	3.9	98	0.00
31		Naphthalene	1.009	0.983	2.6	99	0.00
32		Benzoic acid	0.186	0.216	-16.1	111	0.00
33		4-Chloroaniline	0.449	0.425	5.3	96	-0.01
34	C	Hexachlorobutadiene	0.296	0.310	-4.7	107	-0.01
35		Caprolactam	0.107	0.110	-2.8	101	-0.01
36	C	4-Chloro-3-methylphenol	0.327	0.340	-4.0	105	-0.01
37		2-Methylnaphthalene	0.818	0.793	3.1	99	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.842	0.813	3.4	101	-0.01
40	P	Hexachlorocyclopentadiene	0.444	0.464	-4.5	106	-0.01
41	S	2,4,6-Tribromophenol	0.305	0.321	-5.2	108	-0.02
42	C	2,4,6-Trichlorophenol	0.486	0.491	-1.0	104	-0.01
43		2,4,5-Trichlorophenol	0.538	0.553	-2.8	106	-0.01
44	S	2-Fluorobiphenyl	1.593	1.516	4.8	101	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.708	1.626	4.8	99	-0.01
46		2-Chloronaphthalene	1.303	1.232	5.4	99	-0.01
47		2-Nitroaniline	0.225	0.269	-19.6	122	-0.02
48		Acenaphthylene	2.084	1.986	4.7	100	-0.02
49		Dimethylphthalate	1.762	1.647	6.5	98	-0.01
50		2,6-Dinitrotoluene	0.324	0.361	-11.4	113	-0.01
51	C	Acenaphthene	1.365	1.344	1.5	103	-0.02
52		3-Nitroaniline	0.316	0.342	-8.2	110	-0.01
53	P	2,4-Dinitrophenol	0.115	0.173	-50.4#	159#	-0.01
54		Dibenzofuran	2.093	1.970	5.9	99	-0.02
55	P	4-Nitrophenol	0.257	0.249	3.1	97	-0.01
56		2,4-Dinitrotoluene	0.417	0.497	-19.2	119	-0.01
57		Fluorene	1.700	1.597	6.1	100	-0.01
58		2,3,4,6-Tetrachlorophenol	0.525	0.528	-0.6	105	-0.01
59		Diethylphthalate	1.717	1.582	7.9	97	-0.01
60		4-Chlorophenyl-phenylether	1.040	0.990	4.8	100	-0.02
61		4-Nitroaniline	0.309	0.343	-11.0	108	-0.01
62		Azobenzene	1.101	1.020	7.4	97	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	102	-0.02
64		4,6-Dinitro-2-methylphenol	0.086	0.118	-37.2#	137	-0.01
65	c	n-Nitrosodiphenylamine	0.664	0.628	5.4	97	-0.01
66		4-Bromophenyl-phenylether	0.289	0.282	2.4	100	-0.02
67		Hexachlorobenzene	0.296	0.292	1.4	101	-0.02
68		Atrazine	0.258	0.245	5.0	96	-0.02
69	C	Pentachlorophenol	0.203	0.207	-2.0	99	-0.02
70		Phenanthrene	1.132	1.064	6.0	97	-0.02
71		Anthracene	1.142	1.064	6.8	96	-0.02
72		Carbazole	1.011	0.942	6.8	95	-0.02
73		Di-n-butylphthalate	1.123	1.039	7.5	95	-0.02
74	C	Fluoranthene	1.389	1.298	6.6	97	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	100	-0.04
76		Benzidine	0.617	0.647	-4.9	102	-0.02
77		Pyrene	1.278	1.216	4.9	96	-0.02
78	S	Terphenyl-d14	0.875	0.834	4.7	97	-0.02
79		Butylbenzylphthalate	0.429	0.423	1.4	98	-0.03
80		Benzo(a)anthracene	1.272	1.211	4.8	96	-0.04
81		3,3'-Dichlorobenzidine	0.493	0.472	4.3	95	-0.04
82		Chrysene	1.204	1.134	5.8	95	-0.04
83		Bis(2-ethylhexyl)phthalate	0.621	0.572	7.9	91	-0.04
84	c	Di-n-octyl phthalate	1.046	0.950	9.2	89	-0.06
85		Indeno(1,2,3-cd)pyrene	1.544	1.436	7.0	92	-0.17
86	I	Perylene-d12	1.000	1.000	0.0	96	-0.10
87		Benzo(b)fluoranthene	1.241	1.207	2.7	96	-0.07

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88	Benzo(k)fluoranthene	1.242	1.168	6.0	91	-0.07
89 C	Benzo(a)pyrene	1.189	1.129	5.0	93	-0.09
90	Dibenzo(a,h)anthracene	1.228	1.134	7.7	90	-0.17
91	Benzo(g,h,i)perylene	1.450	1.373	5.3	92	-0.17

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