Evaluate Continuing Calibration Report

Data Path : U:\HPCHEM1\BNA G\Data\BG020518\
Data File : BG032535.D
Acq On : 5 Feb 2018 15:53
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 06 03:32:59 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 29 15:32:55 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	78	-0.01
2		1,4-Dioxane	0.335	0.306	8.7	78	0.00
3		Pyridine	0.905	0.845	6.6	72	0.00
4		n-Nitrosodimethylamine	0.471	0.399	15.3	63	0.00
5	S	2-Fluorophenol	0.996	0.990	0.6	76	0.00
6		Aniline	1.657	1.593	3.9	72	-0.01
7	S	Phenol-d6	1.329	1.227	7.7	72	0.00
8		2-Chlorophenol	1.177	1.197	-1.7	78	-0.01
9		Benzaldehyde	0.686	0.584	14.9	67	-0.01
10	C	Phenol	1.262	1.218	3.5	74	0.00
11		bis(2-Chloroethyl)ether	0.961	0.945	1.7	76	-0.02
12		1,3-Dichlorobenzene	1.592	1.617	-1.6	81	-0.01
13	C	1,4-Dichlorobenzene	1.595	1.581	0.9	78	-0.02
14		1,2-Dichlorobenzene	1.563	1.593	-1.9	79	-0.02
15		Benzyl Alcohol	1.095	1.004	8.3	69	-0.01
16		2,2'-oxybis(1-Chloropropane	0.911	0.861	5.5	73	-0.01
17		2-Methylphenol	1.025	0.933	9.0	70	0.00
18		Hexachloroethane	0.536	0.547	-2.1	82	-0.02
19	P	n-Nitroso-di-n-propylamine	0.881	0.807	8.4	69	-0.02
20		3+4-Methylphenols	1.438	1.370	4.7	70	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	72	-0.01
22		Acetophenone	0.469	0.466	0.6	72	-0.01
23	S	Nitrobenzene-d5	0.320	0.306	4.4	68	-0.01
24		Nitrobenzene	0.310	0.301	2.9	71	-0.01
25		Isophorone	0.545	0.549	-0.7	73	-0.01
26	C	2-Nitrophenol	0.187	0.194	-3.7	71	-0.01
27		2,4-Dimethylphenol	0.269	0.256	4.8	69	0.00
28		bis(2-Chloroethoxy)methane	0.299	0.302	-1.0	73	-0.02
29	C	2,4-Dichlorophenol	0.356	0.334	6.2	68	-0.01
30		1,2,4-Trichlorobenzene	0.474	0.468	1.3	73	-0.02
31		Naphthalene	0.977	1.026	-5.0	77	-0.02
32		Benzoic acid	0.181	0.196	-8.3	76	-0.02
33		4-Chloroaniline	0.436	0.422	3.2	71	-0.01
34	C	Hexachlorobutadiene	0.369	0.365	1.1	73	-0.02
35		Caprolactam	0.102	0.099	2.9	69	0.00
36	C	4-Chloro-3-methylphenol	0.338	0.322	4.7	70	0.00
37		2-Methylnaphthalene	0.821	0.813	1.0	72	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	71	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.936	0.859	8.2	67	-0.02
40	P	Hexachlorocyclopentadiene	0.585	0.511	12.6	62	-0.02
41	S	2,4,6-Tribromophenol	0.381	0.410	-7.6	78	-0.01
42	C	2,4,6-Trichlorophenol	0.509	0.475	6.7	68	-0.02
43		2,4,5-Trichlorophenol	0.593	0.563	5.1	68	0.00
44	S	2-Fluorobiphenyl	1.683	1.587	5.7	69	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.690	1.629	3.6	70	-0.01
46		2-Chloronaphthalene	1.324	1.262	4.7	69	-0.02
47		2-Nitroaniline	0.293	0.266	9.2	65	-0.01
48		Acenaphthylene	2.002	1.935	3.3	70	-0.01
49		Dimethylphthalate	1.844	1.735	5.9	70	-0.01
50		2,6-Dinitrotoluene	0.387	0.365	5.7	70	-0.01
51	C	Acenaphthene	1.389	1.379	0.7	72	-0.01
52		3-Nitroaniline	0.338	0.333	1.5	71	-0.01
53	P	2,4-Dinitrophenol	0.217	0.210	3.2	70	-0.01
54		Dibenzofuran	2.055	1.974	3.9	71	-0.01
55	P	4-Nitrophenol	0.253	0.213	15.8	61	0.00
56		2,4-Dinitrotoluene	0.551	0.544	1.3	71	-0.01
57		Fluorene	1.708	1.680	1.6	74	-0.02
58		2,3,4,6-Tetrachlorophenol	0.585	0.569	2.7	69	-0.02
59		Diethylphthalate	1.796	1.715	4.5	72	-0.02
60		4-Chlorophenyl-phenylether	1.077	1.038	3.6	71	-0.01
61		4-Nitroaniline	0.362	0.351	3.0	73	0.00
62		Azobenzene	1.073	1.050	2.1	73	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	73	-0.01
64		4,6-Dinitro-2-methylphenol	0.149	0.126	15.4	61	-0.01
65	c	n-Nitrosodiphenylamine	0.597	0.587	1.7	73	-0.01
66		4-Bromophenyl-phenylether	0.296	0.291	1.7	72	-0.01
67		Hexachlorobenzene	0.328	0.337	-2.7	77	-0.02
68		Atrazine	0.256	0.252	1.6	71	-0.01
69	C	Pentachlorophenol	0.205	0.194	5.4	70	-0.01
70		Phenanthrene	1.115	1.115	0.0	75	-0.01
71		Anthracene	1.111	1.138	-2.4	75	-0.01
72		Carbazole	0.981	0.995	-1.4	75	0.00
73		Di-n-butylphthalate	1.086	1.153	-6.2	78	-0.02
74	C	Fluoranthene	1.463	1.471	-0.5	76	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	72	-0.01
76		Benzidine	0.393	0.469	-19.3	77	0.00
77		Pyrene	1.227	1.300	-5.9	77	-0.01
78	S	Terphenyl-d14	0.934	0.990	-6.0	78	-0.01
79		Butylbenzylphthalate	0.387	0.463	-19.6	85	-0.01
80		Benzo(a)anthracene	1.240	1.242	-0.2	73	-0.02
81		3,3'-Dichlorobenzidine	0.480	0.492	-2.5	74	-0.01
82		Chrysene	1.197	1.178	1.6	73	-0.01
83		Bis(2-ethylhexyl)phthalate	0.549	0.594	-8.2	77	-0.01
84	c	Di-n-octyl phthalate	0.870	1.030	-18.4	84	-0.02
85		Indeno(1,2,3-cd)pyrene	1.515	1.817	-19.9	87	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	86	-0.01
87		Benzo(b)fluoranthene	1.208	1.101	8.9	80	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.197	1.117	6.7	83	-0.02
89 C	Benzo(a)pyrene	1.152	1.138	1.2	87	-0.02
90	Dibenzo(a,h)anthracene	1.179	1.185	-0.5	88	-0.02
91	Benzo(g,h,i)perylene	1.170	1.151	1.6	86	-0.02

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