Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042116\
Data File : BG021778.D
Acq On : 20 Apr 2016 10:28
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 20 11:03:45 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 13 15:48:13 2016
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	93	-0.01
2		1,4-Dioxane	0.535	0.484	9.5	83	-0.02
3		Pyridine	1.604	1.523	5.0	91	-0.02
4		n-Nitrosodimethylamine	0.795	0.691	13.1	85	-0.02
5	S	2-Fluorophenol	1.201	1.243	-3.5	97	0.00
6		Aniline	2.373	2.344	1.2	95	-0.01
7	S	Phenol-d6	1.794	1.903	-6.1	101	0.00
8		2-Chlorophenol	1.440	1.493	-3.7	98	-0.01
9		Benzaldehyde	1.037	1.137	-9.6	109	-0.02
10	C	Phenol	1.930	1.969	-2.0	96	0.00
11		bis(2-Chloroethyl)ether	1.544	1.481	4.1	92	-0.02
12		1,3-Dichlorobenzene	1.608	1.637	-1.8	95	-0.01
13	C	1,4-Dichlorobenzene	1.637	1.653	-1.0	95	-0.01
14		1,2-Dichlorobenzene	1.570	1.596	-1.7	96	-0.01
15		Benzyl Alcohol	1.371	1.407	-2.6	98	0.00
16		2,2'-oxybis(1-Chloropropane	3.308	2.968	10.3	86	-0.02
17		2-Methylphenol	1.334	1.402	-5.1	101	0.00
18		Hexachloroethane	0.596	0.617	-3.5	98	-0.01
19	P	n-Nitroso-di-n-propylamine	1.394	1.331	4.5	94	-0.01
20		3+4-Methylphenols	1.826	1.914	-4.8	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.557	0.483	13.3	90	-0.01
23	S	Nitrobenzene-d5	0.389	0.384	1.3	102	0.00
24		Nitrobenzene	0.417	0.397	4.8	99	-0.01
25		Isophorone	0.852	0.760	10.8	97	-0.01
26	C	2-Nitrophenol	0.159	0.181	-13.8	119	-0.01
27		2,4-Dimethylphenol	0.361	0.321	11.1	96	0.00
28		bis(2-Chloroethoxy)methane	0.453	0.421	7.1	98	-0.01
29	C	2,4-Dichlorophenol	0.308	0.317	-2.9	106	0.00
30		1,2,4-Trichlorobenzene	0.331	0.335	-1.2	104	-0.01
31		Naphthalene	1.070	1.042	2.6	101	-0.01
32		Benzoic acid	0.192	0.136	29.2#	77	0.02
33		4-Chloroaniline	0.463	0.454	1.9	105	0.00
34	C	Hexachlorobutadiene	0.214	0.217	-1.4	105	-0.01
35		Caprolactam	0.141	0.128	9.2	96	0.00
36	C	4-Chloro-3-methylphenol	0.387	0.401	-3.6	110	0.00
37		2-Methylnaphthalene	0.735	0.747	-1.6	107	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
39		1,2,4,5-Tetrachlorobenzene	0.625	0.629	-0.6	112	-0.01
40	P	Hexachlorocyclopentadiene	0.347	0.282	18.7	87	-0.02
41	S	2,4,6-Tribromophenol	0.216	0.232	-7.4	123	0.00
42	C	2,4,6-Trichlorophenol	0.399	0.425	-6.5	117	0.00
43		2,4,5-Trichlorophenol	0.430	0.442	-2.8	112	0.00
44	S	2-Fluorobiphenyl	1.365	1.357	0.6	111	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.685	1.544	8.4	102	-0.01
46		2-Chloronaphthalene	1.246	1.242	0.3	111	-0.01
47		2-Nitroaniline	0.427	0.418	2.1	111	0.00
48		Acenaphthylene	2.060	2.024	1.7	111	-0.01
49		Dimethylphthalate	1.729	1.624	6.1	110	0.00
50		2,6-Dinitrotoluene	0.330	0.358	-8.5	121	0.00
51	C	Acenaphthene	1.317	1.280	2.8	109	0.00
52		3-Nitroaniline	0.366	0.361	1.4	112	0.00
53	P	2,4-Dinitrophenol	0.106	0.098	7.5	107	0.00
54		Dibenzofuran	1.798	1.777	1.2	112	0.00
55	P	4-Nitrophenol	0.313	0.299	4.5	104	0.01
56		2,4-Dinitrotoluene	0.446	0.484	-8.5	121	0.00
57		Fluorene	1.606	1.563	2.7	112	0.00
58		2,3,4,6-Tetrachlorophenol	0.357	0.371	-3.9	118	0.00
59		Diethylphthalate	1.805	1.658	8.1	107	-0.01
60		4-Chlorophenyl-phenylether	0.791	0.794	-0.4	115	-0.01
61		4-Nitroaniline	0.402	0.381	5.2	103	0.00
62		Azobenzene	1.547	1.443	6.7	107	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	107	-0.01
64		4,6-Dinitro-2-methylphenol	0.077	0.088	-14.3	118	0.00
65	c	n-Nitrosodiphenylamine	0.600	0.615	-2.5	109	0.00
66		4-Bromophenyl-phenylether	0.214	0.229	-7.0	113	-0.01
67		Hexachlorobenzene	0.226	0.239	-5.8	116	0.00
68		Atrazine	0.237	0.235	0.8	100	0.00
69	C	Pentachlorophenol	0.129	0.106	17.8	84	0.00
70		Phenanthrene	1.102	1.088	1.3	106	0.00
71		Anthracene	1.119	1.096	2.1	104	0.00
72		Carbazole	1.044	0.974	6.7	98	0.00
73		Di-n-butylphthalate	1.329	1.241	6.6	94	-0.02
74	C	Fluoranthene	1.316	1.230	6.5	97	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	94	0.00
76		Benzidine	0.623	0.624	-0.2	93	0.00
77		Pyrene	1.153	1.158	-0.4	97	0.00
78	S	Terphenyl-d14	0.802	0.845	-5.4	101	-0.01
79		Butylbenzylphthalate	0.535	0.547	-2.2	95	-0.02
80		Benzo(a)anthracene	1.151	1.147	0.3	94	-0.01
81		3,3'-Dichlorobenzidine	0.911	0.771	15.4	79	0.00
82		Chrysene	1.106	1.107	-0.1	95	-0.01
83		Bis(2-ethylhexyl)phthalate	0.783	0.779	0.5	94	-0.03
84	c	Di-n-octyl phthalate	1.312	1.329	-1.3	94	-0.03
85		Indeno(1,2,3-cd)pyrene	1.315	1.360	-3.4	97	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	94	0.00
87		Benzo(b)fluoranthene	1.160	1.170	-0.9	93	-0.01

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88	Benzo(k)fluoranthene	1.135	1.132	0.3	95	-0.01
89 C	Benzo(a)pyrene	1.124	1.131	-0.6	96	-0.01
90	Dibenzo(a,h)anthracene	1.127	1.171	-3.9	98	-0.02
91	Benzo(g,h,i)perylene	1.106	1.133	-2.4	96	0.00

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