Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF031815\
Data File : BF077828.D
Acq On : 19 Mar 2015 8:32
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 19 09:26:10 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 19 04:21:27 2015
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	111	0.00
2		1,4-Dioxane	0.520	0.480	7.7	102	0.00
3		Pyridine	1.398	1.418	-1.4	115	0.02
4		n-Nitrosodimethylamine	0.609	0.523	14.1	88	0.01
5	S	2-Fluorophenol	1.146	1.173	-2.4	117	0.02
6		Aniline	2.025	1.989	1.8	112	0.00
7	S	Phenol-d6	1.416	1.446	-2.1	114	0.01
8		2-Chlorophenol	1.364	1.387	-1.7	118	0.01
9		Benzaldehyde	0.580	0.689	-18.8	133	0.01
10	C	Phenol	1.673	1.782	-6.5	122	0.02
11		bis(2-Chloroethyl)ether	1.253	1.194	4.7	107	0.00
12		1,3-Dichlorobenzene	1.517	1.492	1.6	113	0.00
13	C	1,4-Dichlorobenzene	1.576	1.595	-1.2	119	0.00
14		1,2-Dichlorobenzene	1.445	1.461	-1.1	118	0.00
15		Benzyl Alcohol	1.105	1.142	-3.3	116	0.01
16		2,2'-oxybis(1-Chloropropane	1.689	1.691	-0.1	114	0.00
17		2-Methylphenol	1.110	1.099	1.0	110	0.01
18		Hexachloroethane	0.517	0.499	3.5	114	0.00
19	P	n-Nitroso-di-n-propylamine	0.979	0.959	2.0	113	0.00
20		3+4-Methylphenols	1.457	1.482	-1.7	117	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
22		Acetophenone	0.443	0.448	-1.1	115	0.00
23	S	Nitrobenzene-d5	0.305	0.303	0.7	115	0.00
24		Nitrobenzene	0.329	0.332	-0.9	119	0.00
25		Isophorone	0.616	0.601	2.4	109	0.01
26	C	2-Nitrophenol	0.161	0.163	-1.2	107	0.01
27		2,4-Dimethylphenol	0.293	0.297	-1.4	113	0.01
28		bis(2-Chloroethoxy)methane	0.374	0.373	0.3	115	0.00
29	C	2,4-Dichlorophenol	0.279	0.274	1.8	109	0.01
30		1,2,4-Trichlorobenzene	0.313	0.301	3.8	109	0.00
31		Naphthalene	1.027	1.009	1.8	112	0.01
32		Benzoic acid	0.141	0.177	-25.5#	128	0.02
33		4-Chloroaniline	0.417	0.409	1.9	108	0.01
34	C	Hexachlorobutadiene	0.177	0.173	2.3	112	0.00
35		Caprolactam	0.082	0.081	1.2	110	0.01
36	C	4-Chloro-3-methylphenol	0.281	0.282	-0.4	115	0.02
37		2-Methylnaphthalene	0.690	0.685	0.7	110	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
39		1,2,4,5-Tetrachlorobenzene	0.597	0.564	5.5	106	0.01
40	P	Hexachlorocyclopentadiene	0.254	0.149	41.3#	60	0.00
41	S	2,4,6-Tribromophenol	0.191	0.190	0.5	109	0.00
42	C	2,4,6-Trichlorophenol	0.413	0.418	-1.2	108	0.01
43		2,4,5-Trichlorophenol	0.446	0.447	-0.2	112	0.01
44	S	2-Fluorobiphenyl	1.361	1.260	7.4	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.599	1.568	1.9	108	0.01
46		2-Chloronaphthalene	1.299	1.247	4.0	109	0.00
47		2-Nitroaniline	0.323	0.336	-4.0	109	0.00
48		Acenaphthylene	2.161	2.162	-0.0	114	0.00
49		Dimethylphthalate	1.483	1.425	3.9	109	0.00
50		2,6-Dinitrotoluene	0.307	0.304	1.0	109	0.00
51	C	Acenaphthene	1.239	1.237	0.2	111	0.01
52		3-Nitroaniline	0.357	0.382	-7.0	116	0.01
53	P	2,4-Dinitrophenol	0.093	0.054	41.9#	64	0.01
54		Dibenzofuran	1.837	1.783	2.9	108	0.01
55	P	4-Nitrophenol	0.265	0.257	3.0	101	0.01
56		2,4-Dinitrotoluene	0.401	0.412	-2.7	114	0.01
57		Fluorene	1.526	1.514	0.8	110	0.01
58		2,3,4,6-Tetrachlorophenol	0.344	0.349	-1.5	112	0.00
59		Diethylphthalate	1.445	1.400	3.1	108	0.01
60		4-Chlorophenyl-phenylether	0.696	0.639	8.2	104	0.01
61		4-Nitroaniline	0.356	0.378	-6.2	117	0.01
62		Azobenzene	1.381	1.528	-10.6	114	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	118	0.00
64		4,6-Dinitro-2-methylphenol	0.078	0.050	35.9#	65	0.01
65	c	n-Nitrosodiphenylamine	0.666	0.641	3.8	111	0.01
66		4-Bromophenyl-phenylether	0.213	0.203	4.7	109	0.01
67		Hexachlorobenzene	0.235	0.220	6.4	109	0.01
68		Atrazine	0.187	0.176	5.9	105	0.00
69	C	Pentachlorophenol	0.104	0.113	-8.7	116	0.01
70		Phenanthrene	1.148	1.126	1.9	115	0.00
71		Anthracene	1.134	1.093	3.6	116	0.00
72		Carbazole	1.079	1.082	-0.3	122	0.01
73		Di-n-butylphthalate	1.210	1.167	3.6	113	0.00
74	C	Fluoranthene	1.266	1.244	1.7	109	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	124	-0.08
76		Benzidine	0.546	0.538	1.5	131	0.00
77		Pyrene	1.258	1.143	9.1	104	0.00
78	S	Terphenyl-d14	0.819	0.709	13.4	102	-0.01
79		Butylbenzylphthalate	0.496	0.488	1.6	120	-0.05
80		Benzo(a)anthracene	1.186	1.128	4.9	123	-0.08
81		3,3'-Dichlorobenzidine	0.391	0.408	-4.3	136	-0.08
82		Chrysene	1.066	1.070	-0.4	126	-0.08
83		Bis(2-ethylhexyl)phthalate	0.751	0.714	4.9	121	-0.09
84	c	Di-n-octyl phthalate	1.284	1.269	1.2	126	-0.17
85		Indeno(1,2,3-cd)pyrene	1.331	1.281	3.8	129	-0.30
86	I	Perylene-d12	1.000	1.000	0.0	129	-0.25
87		Benzo(b)fluoranthene	1.306	1.128	13.6	112	-0.19

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88	Benzo(k)fluoranthene	1.020	1.128	-10.6	151#	-0.19
89 C	Benzo(a)pyrene	1.089	1.084	0.5	130	-0.23
90	Dibenzo(a,h)anthracene	1.081	1.068	1.2	130	-0.31
91	Benzo(g,h,i)perylene	1.100	1.060	3.6	125	-0.31

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