Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF040615\
Data File : BF078271.D
Acq On : 6 Apr 2015 22:02
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 07 02:21:11 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 07 02:08:45 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	118	0.00
2		1,4-Dioxane	0.549	0.503	8.4	108	-0.02
3		Pyridine	1.437	1.481	-3.1	116	-0.03
4		n-Nitrosodimethylamine	0.553	0.601	-8.7	131	-0.03
5	S	2-Fluorophenol	1.213	1.188	2.1	117	-0.01
6		Aniline	2.156	2.110	2.1	118	0.00
7	S	Phenol-d6	1.520	1.509	0.7	120	0.01
8		2-Chlorophenol	1.434	1.395	2.7	116	0.00
9		Benzaldehyde	1.008	0.897	11.0	103	-0.01
10	C	Phenol	1.822	1.765	3.1	115	0.00
11		bis(2-Chloroethyl)ether	1.310	1.320	-0.8	117	0.00
12		1,3-Dichlorobenzene	1.576	1.543	2.1	119	0.00
13	C	1,4-Dichlorobenzene	1.586	1.557	1.8	120	0.00
14		1,2-Dichlorobenzene	1.487	1.439	3.2	117	0.00
15		Benzyl Alcohol	1.068	1.085	-1.6	121	0.00
16		2,2'-oxybis(1-Chloropropane	1.747	1.723	1.4	118	0.00
17		2-Methylphenol	1.114	1.075	3.5	113	0.00
18		Hexachloroethane	0.535	0.546	-2.1	122	-0.01
19	P	n-Nitroso-di-n-propylamine	1.003	0.985	1.8	116	0.00
20		3+4-Methylphenols	1.486	1.475	0.7	116	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	119	0.00
22		Acetophenone	0.466	0.463	0.6	117	-0.01
23	S	Nitrobenzene-d5	0.330	0.319	3.3	114	0.00
24		Nitrobenzene	0.345	0.346	-0.3	121	-0.01
25		Isophorone	0.626	0.636	-1.6	115	0.00
26	C	2-Nitrophenol	0.164	0.165	-0.6	109	0.00
27		2,4-Dimethylphenol	0.306	0.314	-2.6	118	0.00
28		bis(2-Chloroethoxy)methane	0.402	0.404	-0.5	117	0.00
29	C	2,4-Dichlorophenol	0.278	0.282	-1.4	118	0.00
30		1,2,4-Trichlorobenzene	0.319	0.311	2.5	118	-0.01
31		Naphthalene	1.041	1.008	3.2	116	-0.01
32		Benzoic acid	0.132	0.174	-31.8#	149	0.01
33		4-Chloroaniline	0.432	0.446	-3.2	117	0.00
34	C	Hexachlorobutadiene	0.175	0.179	-2.3	121	-0.01
35		Caprolactam	0.083	0.086	-3.6	115	0.01
36	C	4-Chloro-3-methylphenol	0.281	0.288	-2.5	117	0.01
37		2-Methylnaphthalene	0.707	0.674	4.7	113	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	117	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.620	0.608	1.9	113	-0.01
40	P	Hexachlorocyclopentadiene	0.227	0.250	-10.1	124	-0.01
41	S	2,4,6-Tribromophenol	0.166	0.186	-12.0	125	-0.01
42	C	2,4,6-Trichlorophenol	0.391	0.428	-9.5	124	0.00
43		2,4,5-Trichlorophenol	0.408	0.442	-8.3	124	0.00
44	S	2-Fluorobiphenyl	1.399	1.407	-0.6	119	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.636	1.604	2.0	114	-0.01
46		2-Chloronaphthalene	1.287	1.318	-2.4	121	0.00
47		2-Nitroaniline	0.318	0.351	-10.4	123	0.00
48		Acenaphthylene	2.079	2.138	-2.8	120	-0.01
49		Dimethylphthalate	1.503	1.522	-1.3	121	0.00
50		2,6-Dinitrotoluene	0.309	0.322	-4.2	118	-0.01
51	C	Acenaphthene	1.170	1.205	-3.0	123	0.00
52		3-Nitroaniline	0.352	0.399	-13.4	123	0.00
53	P	2,4-Dinitrophenol	0.083	0.073	12.0	99	0.00
54		Dibenzofuran	1.787	1.831	-2.5	122	0.00
55	P	4-Nitrophenol	0.226	0.241	-6.6	113	0.00
56		2,4-Dinitrotoluene	0.400	0.428	-7.0	118	-0.01
57		Fluorene	1.449	1.499	-3.5	120	0.00
58		2,3,4,6-Tetrachlorophenol	0.303	0.333	-9.9	123	0.00
59		Diethylphthalate	1.449	1.495	-3.2	118	-0.01
60		4-Chlorophenyl-phenylether	0.666	0.697	-4.7	123	-0.01
61		4-Nitroaniline	0.355	0.396	-11.5	119	0.00
62		Azobenzene	1.322	1.526	-15.4	136	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	124	-0.01
64		4,6-Dinitro-2-methylphenol	0.087	0.088	-1.1	122	0.00
65	c	n-Nitrosodiphenylamine	0.692	0.695	-0.4	122	-0.01
66		4-Bromophenyl-phenylether	0.212	0.213	-0.5	121	0.00
67		Hexachlorobenzene	0.232	0.225	3.0	118	-0.01
68		Atrazine	0.200	0.205	-2.5	117	-0.01
69	C	Pentachlorophenol	0.085	0.096	-12.9	127	0.00
70		Phenanthrene	1.157	1.160	-0.3	121	-0.01
71		Anthracene	1.140	1.150	-0.9	122	-0.01
72		Carbazole	1.068	1.067	0.1	117	-0.01
73		Di-n-butylphthalate	1.210	1.315	-8.7	126	-0.01
74	C	Fluoranthene	1.220	1.249	-2.4	124	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	124	-0.01
76		Benzidine	0.770	0.625	18.8	99	-0.01
77		Pyrene	1.256	1.245	0.9	124	-0.02
78	S	Terphenyl-d14	0.885	0.854	3.5	119	-0.01
79		Butylbenzylphthalate	0.479	0.529	-10.4	127	-0.01
80		Benzo(a)anthracene	1.190	1.170	1.7	117	-0.01
81		3,3'-Dichlorobenzidine	0.399	0.401	-0.5	121	-0.01
82		Chrysene	1.118	1.101	1.5	126	0.00
83		Bis(2-ethylhexyl)phthalate	0.751	0.789	-5.1	122	-0.01
84	c	Di-n-octyl phthalate	1.232	1.319	-7.1	124	0.00
85		Indeno(1,2,3-cd)pyrene	1.269	1.289	-1.6	124	0.01
86	I	Perylene-d12	1.000	1.000	0.0	122	0.01
87		Benzo(b)fluoranthene	1.236	1.250	-1.1	127	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.098	1.053	4.1	115	0.01
89 C	Benzo(a)pyrene	1.079	1.111	-3.0	123	0.01
90	Dibenzo(a,h)anthracene	1.080	1.110	-2.8	124	0.00
91	Benzo(g,h,i)perylene	1.083	1.109	-2.4	125	0.00

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