Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF031115\
Data File : BF077700.D
Acq On : 11 Mar 2015 21:27
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040EC

Quant Time: Mar 12 05:52:34 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 12 02:24:12 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	107	0.00
2		1,4-Dioxane	0.520	0.474	8.8	97	0.00
3		Pyridine	1.398	1.415	-1.2	112	-0.01
4		n-Nitrosodimethylamine	0.609	0.587	3.6	96	-0.01
5	S	2-Fluorophenol	1.146	1.136	0.9	110	-0.01
6		Aniline	2.025	1.986	1.9	109	-0.01
7	S	Phenol-d6	1.416	1.408	0.6	108	0.00
8		2-Chlorophenol	1.364	1.364	0.0	112	0.00
9		Benzaldehyde	0.580	0.603	-4.0	113	0.00
10	C	Phenol	1.673	1.720	-2.8	114	0.00
11		bis(2-Chloroethyl)ether	1.253	1.254	-0.1	109	0.00
12		1,3-Dichlorobenzene	1.517	1.491	1.7	110	0.00
13	C	1,4-Dichlorobenzene	1.576	1.529	3.0	110	0.00
14		1,2-Dichlorobenzene	1.445	1.377	4.7	108	0.00
15		Benzyl Alcohol	1.105	1.115	-0.9	110	0.00
16		2,2'-oxybis(1-Chloropropane	1.689	1.714	-1.5	112	0.00
17		2-Methylphenol	1.110	1.094	1.4	107	0.00
18		Hexachloroethane	0.517	0.525	-1.5	116	0.00
19	P	n-Nitroso-di-n-propylamine	0.979	0.953	2.7	109	0.00
20		3+4-Methylphenols	1.457	1.458	-0.1	112	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
22		Acetophenone	0.443	0.439	0.9	108	0.00
23	S	Nitrobenzene-d5	0.305	0.305	0.0	111	-0.01
24		Nitrobenzene	0.329	0.322	2.1	111	-0.01
25		Isophorone	0.616	0.613	0.5	107	-0.01
26	C	2-Nitrophenol	0.161	0.164	-1.9	103	0.00
27		2,4-Dimethylphenol	0.293	0.287	2.0	105	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.366	2.1	108	-0.01
29	C	2,4-Dichlorophenol	0.279	0.281	-0.7	107	0.00
30		1,2,4-Trichlorobenzene	0.313	0.310	1.0	107	0.00
31		Naphthalene	1.027	1.017	1.0	108	0.00
32		Benzoic acid	0.141	0.140	0.7	97	-0.01
33		4-Chloroaniline	0.417	0.402	3.6	102	0.00
34	C	Hexachlorobutadiene	0.177	0.175	1.1	109	0.00
35		Caprolactam	0.082	0.080	2.4	105	-0.01
36	C	4-Chloro-3-methylphenol	0.281	0.271	3.6	106	0.00
37		2-Methylnaphthalene	0.690	0.683	1.0	105	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	95	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.597	0.664	-11.2	107	0.00
40	P	Hexachlorocyclopentadiene	0.254	0.300	-18.1	104	0.00
41	S	2,4,6-Tribromophenol	0.191	0.200	-4.7	99	0.00
42	C	2,4,6-Trichlorophenol	0.413	0.456	-10.4	102	0.00
43		2,4,5-Trichlorophenol	0.446	0.478	-7.2	104	0.00
44	S	2-Fluorobiphenyl	1.361	1.461	-7.3	107	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.599	1.810	-13.2	108	0.00
46		2-Chloronaphthalene	1.299	1.412	-8.7	107	-0.01
47		2-Nitroaniline	0.323	0.363	-12.4	102	0.00
48		Acenaphthylene	2.161	2.203	-1.9	100	0.00
49		Dimethylphthalate	1.483	1.499	-1.1	99	0.00
50		2,6-Dinitrotoluene	0.307	0.317	-3.3	99	-0.01
51	C	Acenaphthene	1.239	1.263	-1.9	98	0.00
52		3-Nitroaniline	0.357	0.348	2.5	91	-0.01
53	P	2,4-Dinitrophenol	0.093	0.099	-6.5	101	-0.01
54		Dibenzofuran	1.837	1.806	1.7	95	0.00
55	P	4-Nitrophenol	0.265	0.281	-6.0	95	0.00
56		2,4-Dinitrotoluene	0.401	0.408	-1.7	97	-0.01
57		Fluorene	1.526	1.540	-0.9	96	0.00
58		2,3,4,6-Tetrachlorophenol	0.344	0.361	-4.9	100	0.00
59		Diethylphthalate	1.445	1.456	-0.8	97	0.00
60		4-Chlorophenyl-phenylether	0.696	0.691	0.7	97	0.00
61		4-Nitroaniline	0.356	0.354	0.6	94	-0.01
62		Azobenzene	1.381	1.367	1.0	88	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	-0.01
64		4,6-Dinitro-2-methylphenol	0.078	0.089	-14.1	101	-0.01
65	c	n-Nitrosodiphenylamine	0.666	0.641	3.8	95	0.00
66		4-Bromophenyl-phenylether	0.213	0.211	0.9	97	0.00
67		Hexachlorobenzene	0.235	0.227	3.4	97	0.00
68		Atrazine	0.187	0.189	-1.1	97	0.00
69	C	Pentachlorophenol	0.104	0.114	-9.6	100	0.00
70		Phenanthrene	1.148	1.119	2.5	98	0.00
71		Anthracene	1.134	1.091	3.8	100	-0.01
72		Carbazole	1.079	1.031	4.4	100	-0.01
73		Di-n-butylphthalate	1.210	1.159	4.2	97	-0.01
74	C	Fluoranthene	1.266	1.207	4.7	91	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	98	-0.14
76		Benzidine	0.546	0.470	13.9	91	-0.02
77		Pyrene	1.258	1.243	1.2	90	-0.01
78	S	Terphenyl-d14	0.819	0.781	4.6	89	-0.03
79		Butylbenzylphthalate	0.496	0.489	1.4	95	-0.06
80		Benzo(a)anthracene	1.186	1.167	1.6	100	-0.14
81		3,3'-Dichlorobenzidine	0.391	0.385	1.5	102	-0.14
82		Chrysene	1.066	1.047	1.8	97	-0.14
83		Bis(2-ethylhexyl)phthalate	0.751	0.739	1.6	99	-0.16
84	c	Di-n-octyl phthalate	1.284	1.246	3.0	98	-0.31
85		Indeno(1,2,3-cd)pyrene	1.331	1.286	3.4	102	-0.59#
86	I	Perylene-d12	1.000	1.000	0.0	100	-0.48
87		Benzo(b)fluoranthene	1.306	1.169	10.5	90	-0.37

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88	Benzo(k)fluoranthene	1.020	1.192	-16.9	123	-0.38
89 C	Benzo(a)pyrene	1.089	1.079	0.9	100	-0.46
90	Dibenzo(a,h)anthracene	1.081	1.086	-0.5	102	-0.61#
91	Benzo(g,h,i)perylene	1.100	1.101	-0.1	100	-0.62#

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