Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120315\
Data File : BF083459.D
Acq On : 3 Dec 2015 14:26
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 03 18:21:07 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 03 17:02:53 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	117	-0.01
2		1,4-Dioxane	0.722	0.611	15.4	103	-0.01
3		Pyridine	1.805	1.861	-3.1	131	-0.02
4		n-Nitrosodimethylamine	0.889	0.858	3.5	103	-0.01
5	S	2-Fluorophenol	1.333	1.311	1.7	113	-0.01
6		Aniline	2.496	2.327	6.8	111	-0.02
7	S	Phenol-d6	1.824	1.661	8.9	106	-0.02
8		2-Chlorophenol	1.470	1.348	8.3	105	-0.02
9		Benzaldehyde	0.920	0.834	9.3	110	-0.02
10	C	Phenol	2.195	1.896	13.6	102	-0.01
11		bis(2-Chloroethyl)ether	1.597	1.526	4.4	115	-0.01
12		1,3-Dichlorobenzene	1.634	1.543	5.6	111	-0.02
13	C	1,4-Dichlorobenzene	1.689	1.568	7.2	108	-0.02
14		1,2-Dichlorobenzene	1.550	1.428	7.9	114	-0.02
15		Benzyl Alcohol	1.410	1.322	6.2	107	-0.02
16		2,2'-oxybis(1-Chloropropane	2.802	2.600	7.2	109	-0.02
17		2-Methylphenol	1.225	1.196	2.4	113	-0.01
18		Hexachloroethane	0.587	0.556	5.3	114	-0.02
19	P	n-Nitroso-di-n-propylamine	1.304	1.175	9.9	102	-0.02
20		3+4-Methylphenols	1.660	1.606	3.3	114	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	116	-0.01
22		Acetophenone	0.598	0.540	9.7	105	-0.02
23	S	Nitrobenzene-d5	0.455	0.432	5.1	114	-0.02
24		Nitrobenzene	0.486	0.453	6.8	112	-0.02
25		Isophorone	0.881	0.819	7.0	109	-0.01
26	C	2-Nitrophenol	0.198	0.207	-4.5	126	-0.01
27		2,4-Dimethylphenol	0.332	0.303	8.7	102	-0.02
28		bis(2-Chloroethoxy)methane	0.502	0.498	0.8	121	-0.01
29	C	2,4-Dichlorophenol	0.320	0.306	4.4	116	-0.01
30		1,2,4-Trichlorobenzene	0.348	0.326	6.3	112	-0.02
31		Naphthalene	1.085	1.041	4.1	118	-0.01
32		Benzoic acid	0.229	0.226	1.3	110	0.01
33		4-Chloroaniline	0.468	0.401	14.3	97	-0.02
34	C	Hexachlorobutadiene	0.198	0.187	5.6	116	-0.01
35		Caprolactam	0.101	0.102	-1.0	116	0.00
36	C	4-Chloro-3-methylphenol	0.359	0.329	8.4	111	-0.02
37		2-Methylnaphthalene	0.725	0.675	6.9	112	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.634	0.634	0.0	120	-0.01
40	P	Hexachlorocyclopentadiene	0.301	0.198	34.2#	75	-0.01
41	S	2,4,6-Tribromophenol	0.201	0.189	6.0	111	-0.02
42	C	2,4,6-Trichlorophenol	0.445	0.432	2.9	111	-0.02
43		2,4,5-Trichlorophenol	0.461	0.435	5.6	105	-0.01
44	S	2-Fluorobiphenyl	1.403	1.231	12.3	107	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.754	1.746	0.5	117	-0.02
46		2-Chloronaphthalene	1.336	1.252	6.3	112	-0.02
47		2-Nitroaniline	0.529	0.519	1.9	107	-0.02
48		Acenaphthylene	2.132	1.919	10.0	108	-0.02
49		Dimethylphthalate	1.625	1.510	7.1	119	-0.02
50		2,6-Dinitrotoluene	0.347	0.310	10.7	103	-0.02
51	C	Acenaphthene	1.254	1.110	11.5	107	-0.02
52		3-Nitroaniline	0.407	0.351	13.8	94	-0.02
53	P	2,4-Dinitrophenol	0.187	0.166	11.2	94	-0.01
54		Dibenzofuran	1.838	1.584	13.8	107	-0.02
55	P	4-Nitrophenol	0.254	0.269	-5.9	121	-0.01
56		2,4-Dinitrotoluene	0.440	0.437	0.7	117	-0.01
57		Fluorene	1.451	1.323	8.8	111	-0.02
58		2,3,4,6-Tetrachlorophenol	0.372	0.365	1.9	115	-0.02
59		Diethylphthalate	1.549	1.479	4.5	113	-0.02
60		4-Chlorophenyl-phenylether	0.667	0.647	3.0	113	-0.02
61		4-Nitroaniline	0.407	0.380	6.6	100	-0.02
62		Azobenzene	1.872	1.880	-0.4	114	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	111	-0.03
64		4,6-Dinitro-2-methylphenol	0.134	0.138	-3.0	106	-0.01
65	c	n-Nitrosodiphenylamine	0.703	0.701	0.3	111	-0.02
66		4-Bromophenyl-phenylether	0.219	0.205	6.4	106	-0.03
67		Hexachlorobenzene	0.242	0.224	7.4	105	-0.03
68		Atrazine	0.194	0.176	9.3	112	-0.02
69	C	Pentachlorophenol	0.127	0.121	4.7	104	-0.02
70		Phenanthrene	1.184	1.089	8.0	104	-0.03
71		Anthracene	1.192	1.093	8.3	105	-0.03
72		Carbazole	1.071	1.021	4.7	109	-0.03
73		Di-n-butylphthalate	1.269	1.278	-0.7	117	-0.03
74	C	Fluoranthene	1.227	1.141	7.0	108	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	108	-0.05
76		Benzidine	0.596	0.559	6.2	102	-0.03
77		Pyrene	1.397	1.317	5.7	106	-0.05
78	S	Terphenyl-d14	0.839	0.768	8.5	107	-0.03
79		Butylbenzylphthalate	0.593	0.620	-4.6	120	-0.03
80		Benzo(a)anthracene	1.225	1.164	5.0	106	-0.03
81		3,3'-Dichlorobenzidine	0.383	0.393	-2.6	111	-0.03
82		Chrysene	1.184	1.102	6.9	108	-0.03
83		Bis(2-ethylhexyl)phthalate	0.795	0.869	-9.3	126	-0.05
84	c	Di-n-octyl phthalate	1.184	1.462	-23.5#	133	-0.05
85		Indeno(1,2,3-cd)pyrene	0.862	0.873	-1.3	116	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	122	-0.05
87		Benzo(b)fluoranthene	1.283	1.295	-0.9	129	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.212	1.009	16.7	107	-0.03
89 C	Benzo(a)pyrene	1.102	1.034	6.2	119	-0.05
90	Dibenzo(a,h)anthracene	0.960	0.863	10.1	117	-0.05
91	Benzo(g,h,i)perylene	0.943	0.794	15.8	111	-0.05

(#) = Out of Range SPCC's out = 0 CCC's out = 1