Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090116\
Data File : BF089963.D
Acq On : 1 Sep 2016 10:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 01 13:59:06 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 31 11:55:18 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	110	0.00
2		1,4-Dioxane	0.592	0.577	2.5	113	0.01
3		Pyridine	1.579	1.451	8.1	97	0.01
4		n-Nitrosodimethylamine	0.779	0.770	1.2	105	0.00
5	S	2-Fluorophenol	1.250	1.181	5.5	99	0.00
6		Aniline	2.072	1.991	3.9	111	0.00
7	S	Phenol-d6	1.518	1.453	4.3	104	0.01
8		2-Chlorophenol	1.356	1.353	0.2	108	0.00
9		Benzaldehyde	0.980	0.985	-0.5	110	0.00
10	C	Phenol	1.766	1.650	6.6	105	0.00
11		bis(2-Chloroethyl)ether	1.336	1.398	-4.6	122	0.01
12		1,3-Dichlorobenzene	1.514	1.559	-3.0	116	0.00
13	C	1,4-Dichlorobenzene	1.549	1.528	1.4	113	0.01
14		1,2-Dichlorobenzene	1.389	1.380	0.6	106	0.00
15		Benzyl Alcohol	0.951	0.899	5.5	103	0.00
16		2,2'-oxybis(1-Chloropropane	2.547	2.607	-2.4	110	0.01
17		2-Methylphenol	1.083	1.135	-4.8	115	0.01
18		Hexachloroethane	0.531	0.517	2.6	107	0.00
19	P	n-Nitroso-di-n-propylamine	0.989	0.947	4.2	100	0.00
20		3+4-Methylphenols	1.312	1.299	1.0	123	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	116	0.00
22		Acetophenone	0.449	0.428	4.7	114	0.00
23	S	Nitrobenzene-d5	0.345	0.301	12.8	102	0.00
24		Nitrobenzene	0.367	0.379	-3.3	129	0.01
25		Isophorone	0.701	0.640	8.7	109	0.00
26	C	2-Nitrophenol	0.175	0.167	4.6	111	0.00
27		2,4-Dimethylphenol	0.324	0.323	0.3	127	0.01
28		bis(2-Chloroethoxy)methane	0.423	0.386	8.7	103	0.00
29	C	2,4-Dichlorophenol	0.291	0.287	1.4	121	0.01
30		1,2,4-Trichlorobenzene	0.318	0.306	3.8	109	0.00
31		Naphthalene	0.996	0.948	4.8	120	0.01
32		Benzoic acid	0.218	0.236	-8.3	126	0.01
33		4-Chloroaniline	0.402	0.362	10.0	103	0.00
34	C	Hexachlorobutadiene	0.186	0.178	4.3	115	0.01
35		Caprolactam	0.092	0.097	-5.4	117	0.01
36	C	4-Chloro-3-methylphenol	0.308	0.307	0.3	117	0.00
37		2-Methylnaphthalene	0.658	0.648	1.5	112	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	0.586	0.558	4.8	117	0.01
40	P	Hexachlorocyclopentadiene	0.300	0.334	-11.3	118	0.00
41	S	2,4,6-Tribromophenol	0.197	0.190	3.6	104	0.00
42	C	2,4,6-Trichlorophenol	0.406	0.413	-1.7	105	0.00
43		2,4,5-Trichlorophenol	0.390	0.434	-11.3	131	0.01
44	S	2-Fluorobiphenyl	1.218	1.054	13.5	98	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.610	1.643	-2.0	123	0.01
46		2-Chloronaphthalene	1.139	1.089	4.4	109	0.00
47		2-Nitroaniline	0.372	0.421	-13.2	128	0.01
48		Acenaphthylene	1.884	1.745	7.4	104	0.00
49		Dimethylphthalate	1.446	1.494	-3.3	128	0.01
50		2,6-Dinitrotoluene	0.322	0.335	-4.0	117	0.01
51	C	Acenaphthene	1.193	1.157	3.0	111	0.00
52		3-Nitroaniline	0.367	0.402	-9.5	133	0.01
53	P	2,4-Dinitrophenol	0.130	0.195	-50.0#	148	0.01
54		Dibenzofuran	1.599	1.436	10.2	101	0.00
55	P	4-Nitrophenol	0.277	0.262	5.4	95	0.00
56		2,4-Dinitrotoluene	0.408	0.375	8.1	90	0.00
57		Fluorene	1.169	1.110	5.0	99	0.00
58		2,3,4,6-Tetrachlorophenol	0.334	0.356	-6.6	121	0.01
59		Diethylphthalate	1.360	1.326	2.5	113	0.00
60		4-Chlorophenyl-phenylether	0.550	0.553	-0.5	109	0.00
61		4-Nitroaniline	0.360	0.341	5.3	95	0.00
62		Azobenzene	1.358	1.389	-2.3	120	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
64		4,6-Dinitro-2-methylphenol	0.128	0.149	-16.4	140	0.01
65	c	n-Nitrosodiphenylamine	0.601	0.553	8.0	108	0.00
66		4-Bromophenyl-phenylether	0.201	0.190	5.5	108	0.00
67		Hexachlorobenzene	0.230	0.235	-2.2	126	0.01
68		Atrazine	0.201	0.218	-8.5	130	0.01
69	C	Pentachlorophenol	0.124	0.139	-12.1	111	0.00
70		Phenanthrene	0.996	0.856	14.1	96	0.00
71		Anthracene	1.010	1.036	-2.6	128	0.01
72		Carbazole	0.947	0.843	11.0	103	0.00
73		Di-n-butylphthalate	1.104	1.169	-5.9	125	0.01
74	C	Fluoranthene	1.044	0.917	12.2	95	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	109	0.00
76		Benzidine	0.758	0.728	4.0	105	0.00
77		Pyrene	1.410	1.386	1.7	119	0.01
78	S	Terphenyl-d14	0.810	0.786	3.0	126	0.00
79		Butylbenzylphthalate	0.569	0.587	-3.2	116	0.00
80		Benzo(a)anthracene	1.225	1.147	6.4	105	0.00
81		3,3'-Dichlorobenzidine	0.442	0.512	-15.8	131	0.01
82		Chrysene	1.144	1.080	5.6	110	0.00
83		Bis(2-ethylhexyl)phthalate	0.787	0.831	-5.6	116	0.00
84	c	Di-n-octyl phthalate	1.270	1.442	-13.5	116	0.00
85		Indeno(1,2,3-cd)pyrene	0.846	0.907	-7.2	118	0.01
86	I	Perylene-d12	1.000	1.000	0.0	118	0.01
87		Benzo(b)fluoranthene	1.253	1.090	13.0	91	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.063	1.041	2.1	136	0.00
89 C	Benzo(a)pyrene	1.073	1.001	6.7	108	0.00
90	Dibenzo(a,h)anthracene	0.787	0.777	1.3	116	0.01
91	Benzo(g,h,i)perylene	0.792	0.778	1.8	117	0.01

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