Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110816\
Data File : BF091081.D
Acq On : 8 Nov 2016 13:48
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 09 01:14:03 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 08 12:59:29 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	106	0.00
2		1,4-Dioxane	0.693	0.670	3.3	103	-0.01
3		Pyridine	1.621	1.873	-15.5	116	0.00
4		n-Nitrosodimethylamine	0.641	0.656	-2.3	102	0.00
5	S	2-Fluorophenol	1.211	1.303	-7.6	107	0.00
6		Aniline	1.939	2.022	-4.3	99	0.00
7	S	Phenol-d6	1.644	1.628	1.0	98	0.00
8		2-Chlorophenol	1.445	1.479	-2.4	104	0.00
9		Benzaldehyde	0.912	0.923	-1.2	103	0.00
10	C	Phenol	1.819	1.697	6.7	98	0.00
11		bis(2-Chloroethyl)ether	1.533	1.510	1.5	102	0.00
12		1,3-Dichlorobenzene	1.461	1.509	-3.3	104	0.00
13	C	1,4-Dichlorobenzene	1.568	1.667	-6.3	112	0.00
14		1,2-Dichlorobenzene	1.439	1.390	3.4	95	0.00
15		Benzyl Alcohol	1.159	1.161	-0.2	101	0.00
16		2,2'-oxybis(1-Chloropropane	2.195	2.100	4.3	102	0.00
17		2-Methylphenol	1.144	1.208	-5.6	103	0.00
18		Hexachloroethane	0.607	0.636	-4.8	106	0.00
19	P	n-Nitroso-di-n-propylamine	1.139	1.131	0.7	107	0.00
20		3+4-Methylphenols	1.373	1.555	-13.3	108	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
22		Acetophenone	0.460	0.474	-3.0	103	0.00
23	S	Nitrobenzene-d5	0.367	0.402	-9.5	114	0.00
24		Nitrobenzene	0.405	0.398	1.7	92	0.00
25		Isophorone	0.706	0.711	-0.7	104	0.00
26	C	2-Nitrophenol	0.171	0.200	-17.0	117	0.00
27		2,4-Dimethylphenol	0.283	0.291	-2.8	96	0.01
28		bis(2-Chloroethoxy)methane	0.386	0.384	0.5	91	0.00
29	C	2,4-Dichlorophenol	0.247	0.263	-6.5	104	0.00
30		1,2,4-Trichlorobenzene	0.278	0.294	-5.8	103	0.00
31		Naphthalene	0.956	0.982	-2.7	100	0.00
32		Benzoic acid	0.199	0.218	-9.5	102	0.00
33		4-Chloroaniline	0.398	0.403	-1.3	101	0.00
34	C	Hexachlorobutadiene	0.150	0.170	-13.3	116	0.00
35		Caprolactam	0.078	0.087	-11.5	113	0.00
36	C	4-Chloro-3-methylphenol	0.299	0.293	2.0	94	0.00
37		2-Methylnaphthalene	0.615	0.620	-0.8	104	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
39		1,2,4,5-Tetrachlorobenzene	0.510	0.541	-6.1	97	0.00
40	P	Hexachlorocyclopentadiene	0.164	0.165	-0.6	97	0.00
41	S	2,4,6-Tribromophenol	0.181	0.178	1.7	105	0.00
42	C	2,4,6-Trichlorophenol	0.329	0.345	-4.9	98	0.00
43		2,4,5-Trichlorophenol	0.346	0.369	-6.6	103	0.00
44	S	2-Fluorobiphenyl	1.162	1.262	-8.6	112	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.460	1.459	0.1	98	0.00
46		2-Chloronaphthalene	1.109	1.109	0.0	96	0.00
47		2-Nitroaniline	0.411	0.422	-2.7	93	0.00
48		Acenaphthylene	1.728	1.736	-0.5	101	0.00
49		Dimethylphthalate	1.311	1.235	5.8	96	0.00
50		2,6-Dinitrotoluene	0.281	0.313	-11.4	119	0.00
51	C	Acenaphthene	1.085	1.030	5.1	101	0.00
52		3-Nitroaniline	0.333	0.332	0.3	98	0.00
53	P	2,4-Dinitrophenol	0.132	0.149	-12.9	117	0.00
54		Dibenzofuran	1.460	1.462	-0.1	104	0.00
55	P	4-Nitrophenol	0.186	0.196	-5.4	98	0.00
56		2,4-Dinitrotoluene	0.358	0.400	-11.7	101	0.00
57		Fluorene	1.194	1.191	0.3	100	0.00
58		2,3,4,6-Tetrachlorophenol	0.271	0.302	-11.4	100	0.00
59		Diethylphthalate	1.345	1.317	2.1	96	0.00
60		4-Chlorophenyl-phenylether	0.543	0.590	-8.7	101	0.00
61		4-Nitroaniline	0.337	0.363	-7.7	102	0.00
62		Azobenzene	1.583	1.631	-3.0	106	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	0.122	0.141	-15.6	116	0.00
65	c	n-Nitrosodiphenylamine	0.636	0.633	0.5	105	0.00
66		4-Bromophenyl-phenylether	0.208	0.209	-0.5	109	0.00
67		Hexachlorobenzene	0.229	0.251	-9.6	105	0.00
68		Atrazine	0.183	0.202	-10.4	99	0.00
69	C	Pentachlorophenol	0.102	0.100	2.0	96	0.00
70		Phenanthrene	1.063	1.186	-11.6	105	0.00
71		Anthracene	1.130	1.069	5.4	101	0.00
72		Carbazole	1.019	0.962	5.6	101	0.00
73		Di-n-butylphthalate	1.281	1.328	-3.7	98	0.00
74	C	Fluoranthene	1.103	1.072	2.8	104	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
76		Benzidine	0.447	0.405	9.4	89	0.00
77		Pyrene	1.310	1.269	3.1	113	0.00
78	S	Terphenyl-d14	0.801	0.714	10.9	91	0.00
79		Butylbenzylphthalate	0.627	0.616	1.8	105	0.00
80		Benzo(a)anthracene	1.136	1.129	0.6	107	0.00
81		3,3'-Dichlorobenzidine	0.399	0.368	7.8	99	0.00
82		Chrysene	1.120	0.929	17.1	85	0.00
83		Bis(2-ethylhexyl)phthalate	0.848	0.744	12.3	91	0.00
84	c	Di-n-octyl phthalate	1.394	1.286	7.7	93	0.00
85		Indeno(1,2,3-cd)pyrene	0.929	0.838	9.8	100	0.00
86	I	Perylene-d12	1.000	1.000	0.0	99	0.00
87		Benzo(b)fluoranthene	1.356	1.344	0.9	88	0.00

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88	Benzo(k)fluoranthene	1.190	1.232	-3.5	102	0.00
89 C	Benzo(a)pyrene	1.181	1.184	-0.3	91	0.00
90	Dibenzo(a,h)anthracene	1.021	1.073	-5.1	98	0.00
91	Benzo(g,h,i)perylene	0.923	1.000	-8.3	101	0.01

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