Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120916\
Data File : BF091532.D
Acq On : 9 Dec 2016 17:08
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF120916

Quant Time: Dec 09 19:04:42 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 09 19:00:21 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	101	0.00
2		1,4-Dioxane	0.627	0.600	4.3	101	-0.01
3		Pyridine	1.669	1.689	-1.2	99	-0.01
4		n-Nitrosodimethylamine	0.717	0.673	6.1	96	-0.01
5	S	2-Fluorophenol	1.187	1.209	-1.9	103	0.00
6		Aniline	1.940	1.862	4.0	101	0.00
7	S	Phenol-d6	1.480	1.385	6.4	103	0.00
8		2-Chlorophenol	1.340	1.337	0.2	100	0.00
9		Benzaldehyde	1.018	0.877	13.9	98	0.00
10	C	Phenol	1.730	1.610	6.9	112	0.00
11		bis(2-Chloroethyl)ether	1.300	1.351	-3.9	104	0.00
12		1,3-Dichlorobenzene	1.453	1.335	8.1	101	0.00
13	C	1,4-Dichlorobenzene	1.433	1.367	4.6	101	0.00
14		1,2-Dichlorobenzene	1.292	1.297	-0.4	100	0.00
15		Benzyl Alcohol	1.168	1.054	9.8	96	0.00
16		2,2'-oxybis(1-Chloropropane	2.014	2.005	0.4	99	0.00
17		2-Methylphenol	1.109	1.088	1.9	100	0.00
18		Hexachloroethane	0.559	0.536	4.1	99	0.00
19	P	n-Nitroso-di-n-propylamine	0.925	0.893	3.5	97	-0.01
20		3+4-Methylphenols	1.359	1.134	16.6	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.479	0.460	4.0	103	0.00
23	S	Nitrobenzene-d5	0.358	0.333	7.0	97	0.00
24		Nitrobenzene	0.379	0.398	-5.0	101	0.00
25		Isophorone	0.637	0.619	2.8	101	0.00
26	C	2-Nitrophenol	0.167	0.168	-0.6	97	0.00
27		2,4-Dimethylphenol	0.293	0.269	8.2	101	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.387	-3.5	103	0.00
29	C	2,4-Dichlorophenol	0.255	0.261	-2.4	99	0.00
30		1,2,4-Trichlorobenzene	0.290	0.275	5.2	101	0.00
31		Naphthalene	0.931	0.877	5.8	101	0.00
32		Benzoic acid	0.195	0.205	-5.1	107	0.00
33		4-Chloroaniline	0.401	0.382	4.7	102	0.00
34	C	Hexachlorobutadiene	0.166	0.169	-1.8	102	0.00
35		Caprolactam	0.075	0.074	1.3	109	0.00
36	C	4-Chloro-3-methylphenol	0.286	0.300	-4.9	104	0.00
37		2-Methylnaphthalene	0.602	0.580	3.7	102	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	101	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.549	0.490	10.7	100	0.00
40	P	Hexachlorocyclopentadiene	0.245	0.246	-0.4	100	0.00
41	S	2,4,6-Tribromophenol	0.166	0.151	9.0	95	0.00
42	C	2,4,6-Trichlorophenol	0.352	0.373	-6.0	101	0.00
43		2,4,5-Trichlorophenol	0.362	0.334	7.7	101	0.00
44	S	2-Fluorobiphenyl	1.157	1.096	5.3	100	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.449	1.327	8.4	100	0.00
46		2-Chloronaphthalene	1.122	1.024	8.7	100	0.00
47		2-Nitroaniline	0.377	0.372	1.3	97	-0.01
48		Acenaphthylene	1.710	1.597	6.6	99	0.00
49		Dimethylphthalate	1.266	1.320	-4.3	103	0.00
50		2,6-Dinitrotoluene	0.278	0.300	-7.9	105	0.00
51	C	Acenaphthene	1.188	1.105	7.0	99	0.00
52		3-Nitroaniline	0.327	0.327	0.0	95	0.00
53	P	2,4-Dinitrophenol	0.103	0.123	-19.4	106	0.00
54		Dibenzofuran	1.469	1.366	7.0	98	0.00
55	P	4-Nitrophenol	0.243	0.278	-14.4	100	0.00
56		2,4-Dinitrotoluene	0.348	0.407	-17.0	108	0.00
57		Fluorene	1.127	0.996	11.6	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.284	0.304	-7.0	103	0.00
59		Diethylphthalate	1.219	1.285	-5.4	103	0.00
60		4-Chlorophenyl-phenylether	0.528	0.517	2.1	102	0.00
61		4-Nitroaniline	0.324	0.316	2.5	107	0.00
62		Azobenzene	1.382	1.428	-3.3	101	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
64		4,6-Dinitro-2-methylphenol	0.099	0.105	-6.1	97	-0.01
65	c	n-Nitrosodiphenylamine	0.596	0.543	8.9	103	0.00
66		4-Bromophenyl-phenylether	0.205	0.194	5.4	99	0.00
67		Hexachlorobenzene	0.213	0.200	6.1	104	0.00
68		Atrazine	0.192	0.172	10.4	93	-0.01
69	C	Pentachlorophenol	0.120	0.123	-2.5	102	0.00
70		Phenanthrene	0.990	0.935	5.6	102	0.00
71		Anthracene	1.068	1.019	4.6	101	0.00
72		Carbazole	0.936	0.842	10.0	103	0.00
73		Di-n-butylphthalate	1.091	1.115	-2.2	102	0.00
74	C	Fluoranthene	1.037	1.005	3.1	99	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76		Benzidine	0.618	0.561	9.2	90	-0.01
77		Pyrene	1.587	1.585	0.1	98	0.00
78	S	Terphenyl-d14	0.881	0.880	0.1	98	0.00
79		Butylbenzylphthalate	0.619	0.613	1.0	99	0.00
80		Benzo(a)anthracene	1.162	1.104	5.0	98	0.00
81		3,3'-Dichlorobenzidine	0.417	0.383	8.2	95	0.00
82		Chrysene	1.215	1.142	6.0	94	-0.01
83		Bis(2-ethylhexyl)phthalate	0.794	0.810	-2.0	101	0.00
84	c	Di-n-octyl phthalate	1.347	1.380	-2.4	96	0.00
85		Indeno(1,2,3-cd)pyrene	1.094	1.058	3.3	97	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	96	0.00
87		Benzo(b)fluoranthene	1.192	1.251	-4.9	100	0.00

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88	Benzo(k)fluoranthene	1.048	0.900	14.1	94	0.00
89 C	Benzo(a)pyrene	1.074	1.054	1.9	97	0.00
90	Dibenzo(a,h)anthracene	0.962	0.933	3.0	97	0.00
91	Benzo(g,h,i)perylene	0.979	0.960	1.9	99	0.00

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