Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121816\
Data File : BF091759.D
Acq On : 19 Dec 2016 2:36
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 19 04:06:53 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 17 22:53:54 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	107	0.00
2		1,4-Dioxane	0.588	0.608	-3.4	105	-0.01
3		Pyridine	1.894	1.989	-5.0	101	-0.01
4		n-Nitrosodimethylamine	0.733	0.844	-15.1	115	-0.01
5	S	2-Fluorophenol	1.194	1.242	-4.0	106	0.00
6		Aniline	1.873	1.970	-5.2	108	0.00
7	S	Phenol-d6	1.456	1.496	-2.7	108	0.00
8		2-Chlorophenol	1.314	1.363	-3.7	107	0.00
9		Benzaldehyde	0.903	0.909	-0.7	105	0.00
10	C	Phenol	1.641	1.663	-1.3	104	0.00
11		bis(2-Chloroethyl)ether	1.311	1.306	0.4	107	0.00
12		1,3-Dichlorobenzene	1.443	1.428	1.0	102	-0.01
13	C	1,4-Dichlorobenzene	1.460	1.405	3.8	102	0.00
14		1,2-Dichlorobenzene	1.279	1.403	-9.7	108	0.00
15		Benzyl Alcohol	1.127	1.156	-2.6	103	0.00
16		2,2'-oxybis(1-Chloropropane	2.003	2.106	-5.1	109	0.00
17		2-Methylphenol	1.098	1.128	-2.7	107	0.00
18		Hexachloroethane	0.542	0.571	-5.4	104	0.00
19	P	n-Nitroso-di-n-propylamine	0.968	0.942	2.7	105	0.00
20		3+4-Methylphenols	1.225	1.278	-4.3	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	104	-0.01
22		Acetophenone	0.483	0.509	-5.4	108	0.00
23	S	Nitrobenzene-d5	0.347	0.376	-8.4	102	0.00
24		Nitrobenzene	0.372	0.355	4.6	101	0.00
25		Isophorone	0.662	0.719	-8.6	106	0.00
26	C	2-Nitrophenol	0.180	0.202	-12.2	109	0.00
27		2,4-Dimethylphenol	0.293	0.313	-6.8	106	0.00
28		bis(2-Chloroethoxy)methane	0.387	0.403	-4.1	110	0.00
29	C	2,4-Dichlorophenol	0.270	0.278	-3.0	101	-0.01
30		1,2,4-Trichlorobenzene	0.293	0.302	-3.1	103	-0.01
31		Naphthalene	0.956	0.914	4.4	98	0.00
32		Benzoic acid	0.209	0.213	-1.9	90	0.00
33		4-Chloroaniline	0.402	0.433	-7.7	106	0.00
34	C	Hexachlorobutadiene	0.165	0.176	-6.7	109	0.00
35		Caprolactam	0.087	0.093	-6.9	99	0.00
36	C	4-Chloro-3-methylphenol	0.298	0.315	-5.7	104	0.00
37		2-Methylnaphthalene	0.602	0.665	-10.5	109	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
39		1,2,4,5-Tetrachlorobenzene	0.499	0.501	-0.4	101	-0.01
40	P	Hexachlorocyclopentadiene	0.211	0.194	8.1	83	-0.01
41	S	2,4,6-Tribromophenol	0.170	0.170	0.0	107	0.00
42	C	2,4,6-Trichlorophenol	0.356	0.370	-3.9	102	-0.01
43		2,4,5-Trichlorophenol	0.350	0.357	-2.0	103	0.00
44	S	2-Fluorobiphenyl	0.996	1.056	-6.0	109	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.433	1.528	-6.6	103	0.00
46		2-Chloronaphthalene	1.071	1.158	-8.1	105	0.00
47		2-Nitroaniline	0.366	0.364	0.5	105	0.00
48		Acenaphthylene	1.660	1.773	-6.8	109	0.00
49		Dimethylphthalate	1.306	1.436	-10.0	106	0.00
50		2,6-Dinitrotoluene	0.286	0.329	-15.0	105	0.00
51	C	Acenaphthene	1.139	1.192	-4.7	106	0.00
52		3-Nitroaniline	0.357	0.389	-9.0	110	0.01
53	P	2,4-Dinitrophenol	0.155	0.137	11.6	84	0.00
54		Dibenzofuran	1.355	1.553	-14.6	110	0.00
55	P	4-Nitrophenol	0.248	0.227	8.5	93	0.00
56		2,4-Dinitrotoluene	0.358	0.396	-10.6	103	0.00
57		Fluorene	1.052	1.220	-16.0	107	0.00
58		2,3,4,6-Tetrachlorophenol	0.289	0.285	1.4	101	0.00
59		Diethylphthalate	1.284	1.296	-0.9	103	0.00
60		4-Chlorophenyl-phenylether	0.479	0.469	2.1	101	0.00
61		4-Nitroaniline	0.338	0.347	-2.7	107	0.00
62		Azobenzene	1.366	1.385	-1.4	106	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
64		4,6-Dinitro-2-methylphenol	0.132	0.135	-2.3	88	0.00
65	c	n-Nitrosodiphenylamine	0.621	0.688	-10.8	110	0.00
66		4-Bromophenyl-phenylether	0.210	0.225	-7.1	110	0.00
67		Hexachlorobenzene	0.220	0.242	-10.0	104	0.00
68		Atrazine	0.188	0.197	-4.8	104	0.00
69	C	Pentachlorophenol	0.119	0.127	-6.7	90	0.00
70		Phenanthrene	0.999	1.123	-12.4	102	0.00
71		Anthracene	1.059	1.061	-0.2	112	0.00
72		Carbazole	0.926	1.090	-17.7	113	0.00
73		Di-n-butylphthalate	1.187	1.185	0.2	103	0.00
74	C	Fluoranthene	1.044	1.022	2.1	101	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
76		Benzidine	0.576	0.518	10.1	90	0.00
77		Pyrene	1.444	1.355	6.2	101	0.00
78	S	Terphenyl-d14	0.792	0.738	6.8	94	0.00
79		Butylbenzylphthalate	0.676	0.743	-9.9	115	0.00
80		Benzo(a)anthracene	1.144	1.118	2.3	102	0.00
81		3,3'-Dichlorobenzidine	0.458	0.411	10.3	91	0.00
82		Chrysene	1.186	1.046	11.8	88	-0.01
83		Bis(2-ethylhexyl)phthalate	0.857	0.817	4.7	104	0.00
84	c	Di-n-octyl phthalate	1.580	1.615	-2.2	102	0.00
85		Indeno(1,2,3-cd)pyrene	1.166	1.020	12.5	87	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	89	0.00
87		Benzo(b)fluoranthene	1.268	1.224	3.5	76	0.00

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88	Benzo(k)fluoranthene	0.890	1.216	-36.6#	135	0.00
89 C	Benzo(a)pyrene	1.088	1.133	-4.1	92	0.00
90	Dibenzo(a,h)anthracene	0.949	0.961	-1.3	87	0.00
91	Benzo(g,h,i)perylene	0.964	0.976	-1.2	86	-0.01

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