Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF120717\
Data File : BF101198.D
Acq On : 7 Dec 2017 12:01
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 07 16:32:03 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 04 13:07:49 2017
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	109	0.00
2		1,4-Dioxane	0.500	0.467	6.6	99	0.00
3		Pyridine	1.297	1.217	6.2	92	0.00
4		n-Nitrosodimethylamine	0.634	0.669	-5.5	110	0.00
5	S	2-Fluorophenol	1.210	1.196	1.2	105	0.00
6		Aniline	1.911	1.906	0.3	106	0.00
7	S	Phenol-d6	1.469	1.463	0.4	107	0.00
8		2-Chlorophenol	1.320	1.338	-1.4	109	0.00
9		Benzaldehyde	0.899	0.756	15.9	94	0.00
10	C	Phenol	1.634	1.615	1.2	106	0.00
11		bis(2-Chloroethyl)ether	1.226	1.208	1.5	105	0.00
12		1,3-Dichlorobenzene	1.537	1.540	-0.2	106	0.00
13	C	1,4-Dichlorobenzene	1.591	1.560	1.9	106	0.00
14		1,2-Dichlorobenzene	1.484	1.469	1.0	108	0.00
15		Benzyl Alcohol	1.135	1.136	-0.1	105	0.00
16		2,2'-oxybis(1-Chloropropane	1.666	1.654	0.7	104	0.00
17		2-Methylphenol	1.066	1.060	0.6	107	0.00
18		Hexachloroethane	0.511	0.522	-2.2	110	0.00
19	P	n-Nitroso-di-n-propylamine	0.990	0.987	0.3	109	0.00
20		3+4-Methylphenols	1.390	1.373	1.2	105	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.483	0.483	0.0	107	0.00
23	S	Nitrobenzene-d5	0.335	0.335	0.0	106	0.00
24		Nitrobenzene	0.329	0.328	0.3	107	0.00
25		Isophorone	0.573	0.576	-0.5	108	0.00
26	C	2-Nitrophenol	0.180	0.178	1.1	105	0.00
27		2,4-Dimethylphenol	0.261	0.266	-1.9	109	0.00
28		bis(2-Chloroethoxy)methane	0.385	0.392	-1.8	109	0.00
29	C	2,4-Dichlorophenol	0.284	0.291	-2.5	107	0.00
30		1,2,4-Trichlorobenzene	0.312	0.320	-2.6	110	0.00
31		Naphthalene	0.986	0.981	0.5	107	0.00
32		Benzoic acid	0.203	0.212	-4.4	116	0.01
33		4-Chloroaniline	0.396	0.398	-0.5	105	0.00
34	C	Hexachlorobutadiene	0.192	0.195	-1.6	109	0.00
35		Caprolactam	0.089	0.086	3.4	101	0.00
36	C	4-Chloro-3-methylphenol	0.294	0.299	-1.7	108	0.00
37		2-Methylnaphthalene	0.670	0.664	0.9	105	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
39		1,2,4,5-Tetrachlorobenzene	0.727	0.735	-1.1	108	0.00
40	P	Hexachlorocyclopentadiene	0.215	0.238	-10.7	115	0.00
41	S	2,4,6-Tribromophenol	0.240	0.237	1.3	104	0.00
42	C	2,4,6-Trichlorophenol	0.426	0.430	-0.9	105	0.00
43		2,4,5-Trichlorophenol	0.455	0.471	-3.5	108	0.00
44	S	2-Fluorobiphenyl	1.462	1.458	0.3	106	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.832	1.809	1.3	106	0.00
46		2-Chloronaphthalene	1.301	1.300	0.1	106	0.00
47		2-Nitroaniline	0.385	0.384	0.3	105	0.00
48		Acenaphthylene	2.139	2.087	2.4	104	0.00
49		Dimethylphthalate	1.613	1.605	0.5	106	0.00
50		2,6-Dinitrotoluene	0.340	0.339	0.3	107	0.00
51	C	Acenaphthene	1.281	1.251	2.3	106	0.00
52		3-Nitroaniline	0.382	0.373	2.4	102	0.00
53	P	2,4-Dinitrophenol	0.155	0.148	4.5	96	0.00
54		Dibenzofuran	1.845	1.784	3.3	103	0.00
55	P	4-Nitrophenol	0.258	0.233	9.7	92	0.01
56		2,4-Dinitrotoluene	0.455	0.427	6.2	98	0.00
57		Fluorene	1.500	1.477	1.5	105	0.00
58		2,3,4,6-Tetrachlorophenol	0.365	0.357	2.2	103	0.00
59		Diethylphthalate	1.591	1.578	0.8	104	0.00
60		4-Chlorophenyl-phenylether	0.741	0.742	-0.1	106	0.00
61		4-Nitroaniline	0.397	0.376	5.3	100	0.00
62		Azobenzene	1.350	1.334	1.2	105	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
64		4,6-Dinitro-2-methylphenol	0.134	0.131	2.2	98	0.01
65	c	n-Nitrosodiphenylamine	0.706	0.715	-1.3	104	0.00
66		4-Bromophenyl-phenylether	0.224	0.225	-0.4	103	0.00
67		Hexachlorobenzene	0.240	0.241	-0.4	104	0.00
68		Atrazine	0.216	0.212	1.9	102	0.00
69	C	Pentachlorophenol	0.132	0.128	3.0	95	0.00
70		Phenanthrene	1.101	1.071	2.7	101	0.00
71		Anthracene	1.115	1.093	2.0	100	0.00
72		Carbazole	1.018	0.973	4.4	99	0.00
73		Di-n-butylphthalate	1.277	1.264	1.0	101	0.00
74	C	Fluoranthene	1.221	1.137	6.9	97	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	95	0.00
76		Benzidine	0.650	0.628	3.4	91	0.00
77		Pyrene	1.380	1.448	-4.9	98	0.00
78	S	Terphenyl-d14	0.914	0.959	-4.9	99	0.00
79		Butylbenzylphthalate	0.608	0.645	-6.1	97	0.00
80		Benzo(a)anthracene	1.263	1.268	-0.4	94	0.00
81		3,3'-Dichlorobenzidine	0.437	0.433	0.9	92	0.00
82		Chrysene	1.173	1.183	-0.9	95	0.00
83		Bis(2-ethylhexyl)phthalate	0.868	0.887	-2.2	94	0.00
84	c	Di-n-octyl phthalate	1.444	1.459	-1.0	92	0.00
85		Indeno(1,2,3-cd)pyrene	1.043	0.908	12.9	82	0.03
86	I	Perylene-d12	1.000	1.000	0.0	89	0.01
87		Benzo(b)fluoranthene	1.198	1.227	-2.4	92	0.01

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88	Benzo(k)fluoranthene	1.180	1.194	-1.2	87	0.01
89 C	Benzo(a)pyrene	1.089	1.088	0.1	88	0.01
90	Dibenzo(a,h)anthracene	0.960	0.878	8.5	82	0.02
91	Benzo(g,h,i)perylene	0.905	0.813	10.2	82	0.02

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