Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111116\
Data File : BF091138.D
Acq On : 11 Nov 2016 14:07
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 11 15:53:53 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	104	-0.01
2		1,4-Dioxane	0.693	0.618	10.8	93	-0.05
3		Pyridine	1.621	1.489	8.1	90	-0.03
4		n-Nitrosodimethylamine	0.641	0.543	15.3	83	-0.05
5	S	2-Fluorophenol	1.211	1.326	-9.5	107	-0.01
6		Aniline	1.939	1.903	1.9	92	-0.02
7	S	Phenol-d6	1.644	1.661	-1.0	98	-0.01
8		2-Chlorophenol	1.445	1.496	-3.5	104	-0.02
9		Benzaldehyde	0.912	0.850	6.8	93	-0.01
10	C	Phenol	1.819	1.962	-7.9	111	-0.01
11		bis(2-Chloroethyl)ether	1.533	1.480	3.5	98	-0.02
12		1,3-Dichlorobenzene	1.461	1.511	-3.4	102	-0.02
13	C	1,4-Dichlorobenzene	1.568	1.560	0.5	103	-0.02
14		1,2-Dichlorobenzene	1.439	1.486	-3.3	100	-0.01
15		Benzyl Alcohol	1.159	1.196	-3.2	103	-0.01
16		2,2'-oxybis(1-Chloropropane	2.195	1.866	15.0	89	-0.02
17		2-Methylphenol	1.144	1.106	3.3	92	-0.02
18		Hexachloroethane	0.607	0.585	3.6	95	-0.02
19	P	n-Nitroso-di-n-propylamine	1.139	1.038	8.9	96	-0.02
20		3+4-Methylphenols	1.373	1.441	-5.0	98	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	97	-0.02
22		Acetophenone	0.460	0.490	-6.5	103	-0.01
23	S	Nitrobenzene-d5	0.367	0.354	3.5	96	-0.02
24		Nitrobenzene	0.405	0.422	-4.2	94	-0.01
25		Isophorone	0.706	0.696	1.4	98	-0.01
26	C	2-Nitrophenol	0.171	0.181	-5.8	102	-0.02
27		2,4-Dimethylphenol	0.283	0.309	-9.2	99	-0.01
28		bis(2-Chloroethoxy)methane	0.386	0.429	-11.1	98	-0.01
29	C	2,4-Dichlorophenol	0.247	0.276	-11.7	105	-0.02
30		1,2,4-Trichlorobenzene	0.278	0.307	-10.4	103	-0.01
31		Naphthalene	0.956	0.944	1.3	92	-0.02
32		Benzoic acid	0.199	0.207	-4.0	94	-0.01
33		4-Chloroaniline	0.398	0.416	-4.5	100	-0.01
34	C	Hexachlorobutadiene	0.150	0.179	-19.3	117	-0.02
35		Caprolactam	0.078	0.081	-3.8	101	-0.02
36	C	4-Chloro-3-methylphenol	0.299	0.323	-8.0	100	-0.01
37		2-Methylnaphthalene	0.615	0.680	-10.6	110	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	101	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.510	0.535	-4.9	94	-0.01
40	P	Hexachlorocyclopentadiene	0.164	0.146	11.0	84	-0.02
41	S	2,4,6-Tribromophenol	0.181	0.204	-12.7	119	-0.01
42	C	2,4,6-Trichlorophenol	0.329	0.368	-11.9	103	-0.01
43		2,4,5-Trichlorophenol	0.346	0.368	-6.4	101	-0.01
44	S	2-Fluorobiphenyl	1.162	1.148	1.2	100	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.460	1.496	-2.5	99	-0.01
46		2-Chloronaphthalene	1.109	1.130	-1.9	96	-0.02
47		2-Nitroaniline	0.411	0.451	-9.7	98	-0.01
48		Acenaphthylene	1.728	1.690	2.2	97	-0.02
49		Dimethylphthalate	1.311	1.446	-10.3	111	-0.01
50		2,6-Dinitrotoluene	0.281	0.299	-6.4	112	-0.02
51	C	Acenaphthene	1.085	1.110	-2.3	107	-0.02
52		3-Nitroaniline	0.333	0.382	-14.7	111	-0.01
53	P	2,4-Dinitrophenol	0.132	0.131	0.8	101	-0.01
54		Dibenzofuran	1.460	1.532	-4.9	108	-0.01
55	P	4-Nitrophenol	0.186	0.164	11.8	80	-0.01
56		2,4-Dinitrotoluene	0.358	0.386	-7.8	96	-0.02
57		Fluorene	1.194	1.154	3.4	96	-0.02
58		2,3,4,6-Tetrachlorophenol	0.271	0.286	-5.5	93	-0.01
59		Diethylphthalate	1.345	1.350	-0.4	97	-0.02
60		4-Chlorophenyl-phenylether	0.543	0.608	-12.0	102	-0.02
61		4-Nitroaniline	0.337	0.387	-14.8	107	-0.01
62		Azobenzene	1.583	1.610	-1.7	103	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	103	-0.01
64		4,6-Dinitro-2-methylphenol	0.122	0.121	0.8	102	-0.02
65	c	n-Nitrosodiphenylamine	0.636	0.699	-9.9	119	-0.01
66		4-Bromophenyl-phenylether	0.208	0.200	3.8	107	-0.01
67		Hexachlorobenzene	0.229	0.234	-2.2	100	-0.02
68		Atrazine	0.183	0.161	12.0	80	-0.02
69	C	Pentachlorophenol	0.102	0.096	5.9	94	-0.01
70		Phenanthrene	1.063	1.118	-5.2	101	-0.02
71		Anthracene	1.130	1.100	2.7	106	-0.02
72		Carbazole	1.019	0.995	2.4	107	-0.01
73		Di-n-butylphthalate	1.281	1.243	3.0	93	-0.02
74	C	Fluoranthene	1.103	1.026	7.0	101	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	100	-0.02
76		Benzidine	0.447	0.418	6.5	86	-0.01
77		Pyrene	1.310	1.271	3.0	106	-0.02
78	S	Terphenyl-d14	0.801	0.789	1.5	94	-0.02
79		Butylbenzylphthalate	0.627	0.576	8.1	92	-0.02
80		Benzo(a)anthracene	1.136	1.079	5.0	95	-0.02
81		3,3'-Dichlorobenzidine	0.399	0.395	1.0	99	-0.02
82		Chrysene	1.120	1.050	6.3	90	-0.02
83		Bis(2-ethylhexyl)phthalate	0.848	0.713	15.9	82	-0.02
84	c	Di-n-octyl phthalate	1.394	1.187	14.8	80	-0.02
85		Indeno(1,2,3-cd)pyrene	0.929	0.776	16.5	87	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	93	-0.03
87		Benzo(b)fluoranthene	1.356	1.579	-16.4	98	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.190	1.114	6.4	87	-0.02
89 C	Benzo(a)pyrene	1.181	1.216	-3.0	89	-0.02
90	Dibenzo(a,h)anthracene	1.021	0.988	3.2	86	-0.06
91	Benzo(g,h,i)perylene	0.923	0.926	-0.3	89	-0.05

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