Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092116\
Data File : BF090154.D
Acq On : 21 Sep 2016 11:42
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 22 08:59:47 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 20 17:58:29 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	97	-0.02
2		1,4-Dioxane	0.666	0.590	11.4	94	-0.03
3		Pyridine	1.444	1.396	3.3	98	-0.03
4		n-Nitrosodimethylamine	0.431	0.467	-8.4	111	-0.03
5	S	2-Fluorophenol	1.227	1.234	-0.6	100	-0.01
6		Aniline	1.889	1.892	-0.2	101	-0.01
7	S	Phenol-d6	1.515	1.491	1.6	97	-0.01
8		2-Chlorophenol	1.412	1.352	4.2	97	-0.02
9		Benzaldehyde	0.660	0.678	-2.7	105	-0.01
10	C	Phenol	1.791	1.568	12.5	84	-0.02
11		bis(2-Chloroethyl)ether	1.397	1.275	8.7	90	-0.02
12		1,3-Dichlorobenzene	1.475	1.431	3.0	102	-0.01
13	C	1,4-Dichlorobenzene	1.506	1.448	3.9	102	-0.01
14		1,2-Dichlorobenzene	1.342	1.168	13.0	88	-0.02
15		Benzyl Alcohol	0.904	0.814	10.0	89	-0.02
16		2,2'-oxybis(1-Chloropropane	1.669	1.653	1.0	101	-0.01
17		2-Methylphenol	1.112	1.060	4.7	95	-0.02
18		Hexachloroethane	0.537	0.490	8.8	91	-0.02
19	P	n-Nitroso-di-n-propylamine	0.890	0.908	-2.0	102	-0.01
20		3+4-Methylphenols	1.335	1.298	2.8	97	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	91	-0.02
22		Acetophenone	0.482	0.492	-2.1	97	-0.01
23	S	Nitrobenzene-d5	0.345	0.359	-4.1	103	-0.01
24		Nitrobenzene	0.359	0.329	8.4	85	-0.02
25		Isophorone	0.721	0.704	2.4	97	-0.01
26	C	2-Nitrophenol	0.177	0.179	-1.1	98	-0.01
27		2,4-Dimethylphenol	0.315	0.300	4.8	94	-0.01
28		bis(2-Chloroethoxy)methane	0.436	0.412	5.5	91	-0.02
29	C	2,4-Dichlorophenol	0.276	0.288	-4.3	99	-0.01
30		1,2,4-Trichlorobenzene	0.276	0.276	0.0	95	-0.01
31		Naphthalene	0.952	0.915	3.9	92	-0.01
32		Benzoic acid	0.216	0.223	-3.2	91	-0.02
33		4-Chloroaniline	0.395	0.351	11.1	83	-0.02
34	C	Hexachlorobutadiene	0.145	0.127	12.4	86	-0.02
35		Caprolactam	0.091	0.083	8.8	91	-0.01
36	C	4-Chloro-3-methylphenol	0.312	0.300	3.8	96	-0.01
37		2-Methylnaphthalene	0.592	0.532	10.1	88	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	90	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.459	0.404	12.0	83	-0.02
40	P	Hexachlorocyclopentadiene	0.227	0.199	12.3	75	-0.02
41	S	2,4,6-Tribromophenol	0.172	0.164	4.7	86	-0.02
42	C	2,4,6-Trichlorophenol	0.327	0.307	6.1	81	-0.02
43		2,4,5-Trichlorophenol	0.339	0.341	-0.6	94	-0.01
44	S	2-Fluorobiphenyl	1.019	0.982	3.6	92	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.431	1.254	12.4	78	-0.02
46		2-Chloronaphthalene	1.056	1.009	4.5	95	-0.01
47		2-Nitroaniline	0.318	0.319	-0.3	97	-0.01
48		Acenaphthylene	1.714	1.736	-1.3	98	-0.01
49		Dimethylphthalate	1.274	1.251	1.8	96	-0.01
50		2,6-Dinitrotoluene	0.284	0.267	6.0	83	-0.01
51	C	Acenaphthene	1.018	0.989	2.8	90	-0.01
52		3-Nitroaniline	0.333	0.335	-0.6	94	-0.01
53	P	2,4-Dinitrophenol	0.124	0.127	-2.4	80	-0.02
54		Dibenzofuran	1.339	1.350	-0.8	99	-0.01
55	P	4-Nitrophenol	0.228	0.211	7.5	80	-0.02
56		2,4-Dinitrotoluene	0.336	0.295	12.2	76	-0.02
57		Fluorene	1.009	0.968	4.1	85	-0.02
58		2,3,4,6-Tetrachlorophenol	0.254	0.230	9.4	80	-0.02
59		Diethylphthalate	1.230	1.146	6.8	85	-0.01
60		4-Chlorophenyl-phenylether	0.460	0.423	8.0	90	-0.01
61		4-Nitroaniline	0.339	0.307	9.4	78	-0.02
62		Azobenzene	1.281	1.101	14.1	77	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	85	-0.02
64		4,6-Dinitro-2-methylphenol	0.125	0.129	-3.2	91	-0.01
65	c	n-Nitrosodiphenylamine	0.658	0.616	6.4	85	-0.01
66		4-Bromophenyl-phenylether	0.197	0.200	-1.5	93	-0.01
67		Hexachlorobenzene	0.229	0.233	-1.7	99	-0.01
68		Atrazine	0.180	0.192	-6.7	104	-0.01
69	C	Pentachlorophenol	0.130	0.134	-3.1	87	-0.01
70		Phenanthrene	0.935	0.891	4.7	80	-0.02
71		Anthracene	0.981	1.038	-5.8	97	-0.01
72		Carbazole	1.037	1.062	-2.4	96	-0.01
73		Di-n-butylphthalate	1.206	1.294	-7.3	96	-0.01
74	C	Fluoranthene	1.004	1.055	-5.1	96	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	96	-0.01
76		Benzidine	0.669	0.571	14.6	81	-0.01
77		Pyrene	1.369	1.296	5.3	98	-0.02
78	S	Terphenyl-d14	0.788	0.749	4.9	104	-0.01
79		Butylbenzylphthalate	0.669	0.693	-3.6	103	-0.01
80		Benzo(a)anthracene	1.076	0.990	8.0	88	-0.01
81		3,3'-Dichlorobenzidine	0.444	0.424	4.5	95	-0.01
82		Chrysene	1.048	0.907	13.5	84	-0.02
83		Bis(2-ethylhexyl)phthalate	0.857	0.797	7.0	88	-0.02
84	c	Di-n-octyl phthalate	1.476	1.393	5.6	86	-0.02
85		Indeno(1,2,3-cd)pyrene	0.847	0.908	-7.2	99	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	85	-0.02
87		Benzo(b)fluoranthene	1.107	0.986	10.9	75	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.039	1.010	2.8	91	-0.02
89 C	Benzo(a)pyrene	1.042	0.978	6.1	83	-0.02
90	Dibenzo(a,h)anthracene	0.813	0.894	-10.0	98	-0.03
91	Benzo(g,h,i)perylene	0.796	0.851	-6.9	96	-0.03

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