Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF033116\
Data File : BF085947.D
Acq On : 1 Apr 2016 00:39
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 01 01:14:50 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 30 12:31:18 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	138	-0.01
2		1,4-Dioxane	0.573	0.560	2.3	136	-0.01
3		Pyridine	1.375	1.381	-0.4	132	-0.01
4		n-Nitrosodimethylamine	0.550	0.573	-4.2	138	-0.01
5	S	2-Fluorophenol	1.201	1.145	4.7	138	-0.01
6		Aniline	2.055	1.902	7.4	134	0.00
7	S	Phenol-d6	1.527	1.464	4.1	140	-0.01
8		2-Chlorophenol	1.395	1.316	5.7	130	-0.01
9		Benzaldehyde	0.920	0.809	12.1	127	0.00
10	C	Phenol	1.730	1.535	11.3	128	0.00
11		bis(2-Chloroethyl)ether	1.331	1.288	3.2	139	-0.01
12		1,3-Dichlorobenzene	1.503	1.512	-0.6	138	-0.01
13	C	1,4-Dichlorobenzene	1.524	1.500	1.6	145	0.00
14		1,2-Dichlorobenzene	1.420	1.330	6.3	130	-0.01
15		Benzyl Alcohol	1.019	0.889	12.8	117	-0.01
16		2,2'-oxybis(1-Chloropropane	1.766	1.541	12.7	123	-0.01
17		2-Methylphenol	1.117	1.112	0.4	135	0.00
18		Hexachloroethane	0.558	0.542	2.9	138	-0.01
19	P	n-Nitroso-di-n-propylamine	0.914	0.845	7.5	123	-0.01
20		3+4-Methylphenols	1.343	1.281	4.6	137	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	131	-0.01
22		Acetophenone	0.466	0.459	1.5	139	0.00
23	S	Nitrobenzene-d5	0.355	0.346	2.5	130	-0.01
24		Nitrobenzene	0.366	0.387	-5.7	152#	-0.01
25		Isophorone	0.684	0.664	2.9	136	-0.01
26	C	2-Nitrophenol	0.183	0.185	-1.1	140	0.00
27		2,4-Dimethylphenol	0.320	0.293	8.4	117	-0.01
28		bis(2-Chloroethoxy)methane	0.390	0.361	7.4	121	-0.01
29	C	2,4-Dichlorophenol	0.269	0.273	-1.5	140	-0.01
30		1,2,4-Trichlorobenzene	0.299	0.304	-1.7	140	-0.01
31		Naphthalene	1.008	0.997	1.1	141	0.00
32		Benzoic acid	0.213	0.236	-10.8	134	0.01
33		4-Chloroaniline	0.411	0.391	4.9	130	0.00
34	C	Hexachlorobutadiene	0.162	0.166	-2.5	140	-0.01
35		Caprolactam	0.095	0.085	10.5	129	0.00
36	C	4-Chloro-3-methylphenol	0.314	0.300	4.5	138	0.00
37		2-Methylnaphthalene	0.635	0.558	12.1	122	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	119	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.593	0.673	-13.5	134	-0.01
40	P	Hexachlorocyclopentadiene	0.196	0.110	43.9#	66	-0.01
41	S	2,4,6-Tribromophenol	0.190	0.182	4.2	113	0.00
42	C	2,4,6-Trichlorophenol	0.401	0.421	-5.0	128	-0.01
43		2,4,5-Trichlorophenol	0.420	0.429	-2.1	123	-0.01
44	S	2-Fluorobiphenyl	1.197	1.156	3.4	112	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.688	1.752	-3.8	124	-0.01
46		2-Chloronaphthalene	1.241	1.297	-4.5	120	-0.01
47		2-Nitroaniline	0.379	0.413	-9.0	140	-0.01
48		Acenaphthylene	2.022	1.898	6.1	110	-0.01
49		Dimethylphthalate	1.517	1.655	-9.1	140	0.00
50		2,6-Dinitrotoluene	0.327	0.307	6.1	108	-0.01
51	C	Acenaphthene	1.235	1.121	9.2	116	-0.01
52		3-Nitroaniline	0.374	0.345	7.8	119	-0.01
53	P	2,4-Dinitrophenol	0.140	0.129	7.9	100	-0.01
54		Dibenzofuran	1.599	1.435	10.3	108	-0.01
55	P	4-Nitrophenol	0.244	0.211	13.5	99	-0.01
56		2,4-Dinitrotoluene	0.415	0.423	-1.9	112	0.00
57		Fluorene	1.329	1.230	7.4	108	-0.01
58		2,3,4,6-Tetrachlorophenol	0.292	0.296	-1.4	124	-0.01
59		Diethylphthalate	1.499	1.499	0.0	124	-0.01
60		4-Chlorophenyl-phenylether	0.650	0.615	5.4	124	-0.01
61		4-Nitroaniline	0.358	0.354	1.1	112	-0.01
62		Azobenzene	1.441	1.427	1.0	121	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	106	-0.01
64		4,6-Dinitro-2-methylphenol	0.117	0.114	2.6	98	0.00
65	c	n-Nitrosodiphenylamine	0.633	0.633	0.0	108	-0.01
66		4-Bromophenyl-phenylether	0.205	0.213	-3.9	113	-0.01
67		Hexachlorobenzene	0.214	0.216	-0.9	105	-0.01
68		Atrazine	0.214	0.212	0.9	110	0.00
69	C	Pentachlorophenol	0.105	0.095	9.5	90	0.00
70		Phenanthrene	1.027	1.003	2.3	100	-0.01
71		Anthracene	1.078	1.127	-4.5	120	-0.01
72		Carbazole	0.905	0.888	1.9	102	-0.01
73		Di-n-butylphthalate	1.242	1.158	6.8	105	-0.01
74	C	Fluoranthene	1.092	0.975	10.7	101	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	94	-0.01
76		Benzidine	0.704	0.682	3.1	89	-0.01
77		Pyrene	1.644	1.590	3.3	101	-0.01
78	S	Terphenyl-d14	0.972	0.896	7.8	97	-0.01
79		Butylbenzylphthalate	0.688	0.763	-10.9	107	-0.01
80		Benzo(a)anthracene	1.140	1.205	-5.7	103	-0.01
81		3,3'-Dichlorobenzidine	0.791	0.997	-26.0#	130	-0.01
82		Chrysene	1.149	1.175	-2.3	102	-0.01
83		Bis(2-ethylhexyl)phthalate	0.824	0.951	-15.4	109	-0.01
84	c	Di-n-octyl phthalate	1.226	1.541	-25.7#	116	-0.01
85		Indeno(1,2,3-cd)pyrene	0.887	1.427	-60.9#	153#	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	140	-0.01
87		Benzo(b)fluoranthene	1.260	1.053	16.4	115	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.136	1.148	-1.1	153#	-0.01
89 C	Benzo(a)pyrene	1.119	1.093	2.3	139	-0.01
90	Dibenzo(a,h)anthracene	1.039	1.118	-7.6	155#	-0.01
91	Benzo(g,h,i)perylene	1.052	1.064	-1.1	145	-0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 1