Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF040716\
Data File : BF086125.D
Acq On : 7 Apr 2016 11:55
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 08 01:53:47 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 01 23:17:06 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	109	-0.03
2		1,4-Dioxane	0.555	0.625	-12.6	129	-0.06
3		Pyridine	1.242	1.619	-30.4#	129	-0.08
4		n-Nitrosodimethylamine	0.546	0.582	-6.6	115	-0.07
5	S	2-Fluorophenol	1.170	1.222	-4.4	114	-0.05
6		Aniline	1.950	1.992	-2.2	110	-0.03
7	S	Phenol-d6	1.486	1.504	-1.2	108	-0.03
8		2-Chlorophenol	1.395	1.438	-3.1	116	-0.03
9		Benzaldehyde	0.800	0.832	-4.0	113	-0.05
10	C	Phenol	1.695	1.662	1.9	100	-0.02
11		bis(2-Chloroethyl)ether	1.342	1.318	1.8	120	-0.03
12		1,3-Dichlorobenzene	1.479	1.532	-3.6	116	-0.03
13	C	1,4-Dichlorobenzene	1.524	1.533	-0.6	115	-0.05
14		1,2-Dichlorobenzene	1.402	1.470	-4.9	113	-0.03
15		Benzyl Alcohol	0.962	0.977	-1.6	115	-0.03
16		2,2'-oxybis(1-Chloropropane	1.640	1.674	-2.1	113	-0.03
17		2-Methylphenol	1.100	1.161	-5.5	122	-0.02
18		Hexachloroethane	0.560	0.567	-1.2	115	-0.05
19	P	n-Nitroso-di-n-propylamine	0.921	0.856	7.1	111	-0.03
20		3+4-Methylphenols	1.338	1.479	-10.5	125	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	112	-0.03
22		Acetophenone	0.470	0.485	-3.2	121	-0.03
23	S	Nitrobenzene-d5	0.339	0.331	2.4	117	-0.03
24		Nitrobenzene	0.371	0.410	-10.5	126	-0.03
25		Isophorone	0.669	0.673	-0.6	116	-0.03
26	C	2-Nitrophenol	0.178	0.189	-6.2	124	-0.02
27		2,4-Dimethylphenol	0.319	0.307	3.8	120	-0.02
28		bis(2-Chloroethoxy)methane	0.393	0.370	5.9	117	-0.03
29	C	2,4-Dichlorophenol	0.270	0.274	-1.5	119	-0.03
30		1,2,4-Trichlorobenzene	0.303	0.305	-0.7	118	-0.03
31		Naphthalene	1.006	0.967	3.9	113	-0.03
32		Benzoic acid	0.234	0.161	31.2#	76	-0.03
33		4-Chloroaniline	0.406	0.399	1.7	117	-0.02
34	C	Hexachlorobutadiene	0.163	0.160	1.8	114	-0.03
35		Caprolactam	0.086	0.086	0.0	106	-0.02
36	C	4-Chloro-3-methylphenol	0.303	0.321	-5.9	119	-0.02
37		2-Methylnaphthalene	0.657	0.615	6.4	113	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	114	-0.03
39		1,2,4,5-Tetrachlorobenzene	0.601	0.614	-2.2	118	-0.03
40	P	Hexachlorocyclopentadiene	0.163	0.126	22.7	87	-0.03
41	S	2,4,6-Tribromophenol	0.188	0.177	5.9	106	-0.03
42	C	2,4,6-Trichlorophenol	0.386	0.404	-4.7	120	-0.03
43		2,4,5-Trichlorophenol	0.392	0.407	-3.8	121	-0.03
44	S	2-Fluorobiphenyl	1.203	1.212	-0.7	112	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.724	1.709	0.9	116	-0.03
46		2-Chloronaphthalene	1.250	1.297	-3.8	117	-0.03
47		2-Nitroaniline	0.366	0.411	-12.3	124	-0.03
48		Acenaphthylene	2.002	2.021	-0.9	111	-0.03
49		Dimethylphthalate	1.482	1.479	0.2	123	-0.02
50		2,6-Dinitrotoluene	0.314	0.286	8.9	97	-0.03
51	C	Acenaphthene	1.191	1.179	1.0	111	-0.03
52		3-Nitroaniline	0.378	0.358	5.3	100	-0.03
53	P	2,4-Dinitrophenol	0.127	0.130	-2.4	96	-0.03
54		Dibenzofuran	1.573	1.513	3.8	108	-0.03
55	P	4-Nitrophenol	0.223	0.216	3.1	112	-0.02
56		2,4-Dinitrotoluene	0.396	0.404	-2.0	123	-0.02
57		Fluorene	1.344	1.362	-1.3	112	-0.03
58		2,3,4,6-Tetrachlorophenol	0.292	0.305	-4.5	121	-0.03
59		Diethylphthalate	1.496	1.313	12.2	108	-0.03
60		4-Chlorophenyl-phenylether	0.626	0.591	5.6	109	-0.03
61		4-Nitroaniline	0.372	0.399	-7.3	123	-0.02
62		Azobenzene	1.433	1.421	0.8	114	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	112	-0.03
64		4,6-Dinitro-2-methylphenol	0.121	0.115	5.0	109	-0.02
65	c	n-Nitrosodiphenylamine	0.625	0.617	1.3	110	-0.03
66		4-Bromophenyl-phenylether	0.204	0.211	-3.4	114	-0.03
67		Hexachlorobenzene	0.211	0.225	-6.6	118	-0.03
68		Atrazine	0.202	0.184	8.9	110	-0.03
69	C	Pentachlorophenol	0.099	0.088	11.1	96	-0.03
70		Phenanthrene	1.091	1.091	0.0	112	-0.03
71		Anthracene	1.143	1.035	9.4	114	-0.03
72		Carbazole	0.982	0.882	10.2	109	-0.03
73		Di-n-butylphthalate	1.217	1.175	3.5	107	-0.03
74	C	Fluoranthene	1.135	1.114	1.9	120	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	107	-0.02
76		Benzidine	0.706	0.694	1.7	104	-0.02
77		Pyrene	1.409	1.475	-4.7	117	-0.02
78	S	Terphenyl-d14	0.851	0.813	4.5	110	-0.03
79		Butylbenzylphthalate	0.635	0.717	-12.9	121	-0.03
80		Benzo(a)anthracene	1.179	1.199	-1.7	113	-0.02
81		3,3'-Dichlorobenzidine	0.861	0.940	-9.2	110	-0.03
82		Chrysene	1.136	1.164	-2.5	106	-0.03
83		Bis(2-ethylhexyl)phthalate	0.842	0.897	-6.5	114	-0.02
84	c	Di-n-octyl phthalate	1.345	1.412	-5.0	105	-0.03
85		Indeno(1,2,3-cd)pyrene	0.921	0.852	7.5	100	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	95	-0.03
87		Benzo(b)fluoranthene	1.258	1.492	-18.6	100	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.114	1.001	10.1	103	-0.03
89 C	Benzo(a)pyrene	1.115	1.105	0.9	98	-0.03
90	Dibenzo(a,h)anthracene	0.897	0.905	-0.9	99	-0.06
91	Benzo(g,h,i)perylene	0.868	0.890	-2.5	102	-0.06

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