Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF123114\
Data File : BF076697.D
Acq On : 31 Dec 2014 12:05
Operator : IZ / TP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 01 05:06:43 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121714.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 01 04:58:19 2015
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	99	-0.20
2		1,4-Dioxane	0.504	0.440	12.7	86	-0.27
3		Pyridine	1.269	1.115	12.1	89	-0.39
4		n-Nitrosodimethylamine	0.614	0.685	-11.6	109	-0.37
5	S	2-Fluorophenol	1.180	1.190	-0.8	101	-0.27
6		Aniline	2.058	2.097	-1.9	100	-0.20
7	S	Phenol-d6	1.441	1.427	1.0	97	-0.18
8		2-Chlorophenol	1.351	1.426	-5.6	103	-0.20
9		Benzaldehyde	1.096	0.989	9.8	88	-0.22
10	C	Phenol	1.749	1.699	2.9	95	-0.16
11		bis(2-Chloroethyl)ether	1.259	1.270	-0.9	100	-0.19
12		1,3-Dichlorobenzene	1.568	1.608	-2.6	100	-0.20
13	C	1,4-Dichlorobenzene	1.589	1.604	-0.9	102	-0.20
14		1,2-Dichlorobenzene	1.506	1.589	-5.5	106	-0.20
15		Benzyl Alcohol	1.248	1.294	-3.7	99	-0.19
16		2,2'-oxybis(1-Chloropropane	1.823	1.847	-1.3	99	-0.18
17		2-Methylphenol	1.110	1.099	1.0	95	-0.16
18		Hexachloroethane	0.568	0.595	-4.8	105	-0.20
19	P	n-Nitroso-di-n-propylamine	1.052	1.066	-1.3	104	-0.16
20		3+4-Methylphenols	1.499	1.541	-2.8	100	-0.15
21	I	Naphthalene-d8	1.000	1.000	0.0	98	-0.19
22		Acetophenone	0.479	0.490	-2.3	100	-0.19
23	S	Nitrobenzene-d5	0.336	0.339	-0.9	99	-0.19
24		Nitrobenzene	0.349	0.350	-0.3	98	-0.19
25		Isophorone	0.648	0.665	-2.6	101	-0.19
26	C	2-Nitrophenol	0.170	0.175	-2.9	97	-0.19
27		2,4-Dimethylphenol	0.276	0.290	-5.1	102	-0.16
28		bis(2-Chloroethoxy)methane	0.380	0.392	-3.2	97	-0.18
29	C	2,4-Dichlorophenol	0.271	0.281	-3.7	103	-0.19
30		1,2,4-Trichlorobenzene	0.292	0.291	0.3	97	-0.20
31		Naphthalene	0.991	1.030	-3.9	103	-0.19
32		Benzoic acid	0.153	0.151	1.3	96	-0.15
33		4-Chloroaniline	0.404	0.421	-4.2	99	-0.19
34	C	Hexachlorobutadiene	0.170	0.176	-3.5	105	-0.19
35		Caprolactam	0.078	0.078	0.0	100	-0.19
36	C	4-Chloro-3-methylphenol	0.301	0.321	-6.6	100	-0.18
37		2-Methylnaphthalene	0.691	0.725	-4.9	102	-0.20
38	I	Acenaphthene-d10	1.000	1.000	0.0	101	-0.22
39		1,2,4,5-Tetrachlorobenzene	0.492	0.501	-1.8	102	-0.20
40	P	Hexachlorocyclopentadiene	0.241	0.261	-8.3	103	-0.20
41	S	2,4,6-Tribromophenol	0.169	0.172	-1.8	101	-0.23
42	C	2,4,6-Trichlorophenol	0.357	0.357	0.0	98	-0.20
43		2,4,5-Trichlorophenol	0.384	0.384	0.0	98	-0.20
44	S	2-Fluorobiphenyl	1.285	1.324	-3.0	105	-0.20

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.486	1.485	0.1	97	-0.20
46		2-Chloronaphthalene	1.178	1.149	2.5	99	-0.21
47		2-Nitroaniline	0.358	0.365	-2.0	98	-0.20
48		Acenaphthylene	2.004	2.003	0.0	101	-0.22
49		Dimethylphthalate	1.415	1.432	-1.2	104	-0.20
50		2,6-Dinitrotoluene	0.290	0.298	-2.8	101	-0.20
51	C	Acenaphthene	1.134	1.170	-3.2	105	-0.22
52		3-Nitroaniline	0.329	0.319	3.0	90	-0.21
53	P	2,4-Dinitrophenol	0.128	0.091	28.9#	73	-0.20
54		Dibenzofuran	1.640	1.676	-2.2	103	-0.21
55	P	4-Nitrophenol	0.252	0.237	6.0	102	-0.18
56		2,4-Dinitrotoluene	0.386	0.365	5.4	89	-0.21
57		Fluorene	1.378	1.406	-2.0	103	-0.22
58		2,3,4,6-Tetrachlorophenol	0.286	0.300	-4.9	102	-0.21
59		Diethylphthalate	1.445	1.505	-4.2	105	-0.20
60		4-Chlorophenyl-phenylether	0.596	0.571	4.2	98	-0.22
61		4-Nitroaniline	0.346	0.318	8.1	85	-0.22
62		Azobenzene	1.424	1.553	-9.1	112	-0.22
63	I	Phenanthrene-d10	1.000	1.000	0.0	95	-0.23
64		4,6-Dinitro-2-methylphenol	0.125	0.109	12.8	84	-0.21
65	c	n-Nitrosodiphenylamine	0.703	0.721	-2.6	97	-0.22
66		4-Bromophenyl-phenylether	0.195	0.220	-12.8	106	-0.22
67		Hexachlorobenzene	0.227	0.235	-3.5	98	-0.23
68		Atrazine	0.210	0.216	-2.9	95	-0.21
69	C	Pentachlorophenol	0.111	0.126	-13.5	105	-0.23
70		Phenanthrene	1.207	1.231	-2.0	95	-0.24
71		Anthracene	1.184	1.239	-4.6	102	-0.23
72		Carbazole	1.148	1.185	-3.2	96	-0.23
73		Di-n-butylphthalate	1.410	1.672	-18.6	111	-0.21
74	C	Fluoranthene	1.232	1.232	0.0	97	-0.26
75	I	Chrysene-d12	1.000	1.000	0.0	88	-0.28
76		Benzidine	0.833	0.732	12.1	75	-0.24
77		Pyrene	1.399	1.457	-4.1	94	-0.27
78	S	Terphenyl-d14	0.907	0.998	-10.0	99	-0.24
79		Butylbenzylphthalate	0.669	0.777	-16.1	103	-0.24
80		Benzo(a)anthracene	1.245	1.287	-3.4	93	-0.28
81		3,3'-Dichlorobenzidine	0.416	0.435	-4.6	90	-0.27
82		Chrysene	1.140	1.154	-1.2	90	-0.28
83		Bis(2-ethylhexyl)phthalate	1.012	1.201	-18.7	105	-0.24
84	c	Di-n-octyl phthalate	1.728	2.042	-18.2	106	-0.26
85		Indeno(1,2,3-cd)pyrene	1.251	1.268	-1.4	95	-0.46
86	I	Perylene-d12	1.000	1.000	0.0	95	-0.31
87		Benzo(b)fluoranthene	1.319	1.516	-14.9	108	-0.31

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.229	0.993	19.2	73	-0.31
89 C	Benzo(a)pyrene	1.177	1.161	1.4	91	-0.32
90	Dibenzo(a,h)anthracene	1.167	1.170	-0.3	94	-0.45
91	Benzo(g,h,i)perylene	1.135	1.139	-0.4	94	-0.52#

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