Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF120215\
Data File : BF083437.D
Acq On : 3 Dec 2015 1:37
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 03 06:34:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 01 01:17:16 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	106	0.00
2		1,4-Dioxane	0.722	0.627	13.2	96	-0.01
3		Pyridine	1.805	1.724	4.5	110	-0.02
4		n-Nitrosodimethylamine	0.889	0.854	3.9	93	-0.01
5	S	2-Fluorophenol	1.333	1.311	1.7	102	-0.01
6		Aniline	2.496	2.382	4.6	103	-0.01
7	S	Phenol-d6	1.824	1.683	7.7	97	-0.01
8		2-Chlorophenol	1.470	1.396	5.0	98	-0.01
9		Benzaldehyde	0.920	0.879	4.5	104	-0.01
10	C	Phenol	2.195	2.030	7.5	99	-0.01
11		bis(2-Chloroethyl)ether	1.597	1.551	2.9	105	0.00
12		1,3-Dichlorobenzene	1.634	1.525	6.7	99	-0.01
13	C	1,4-Dichlorobenzene	1.689	1.601	5.2	99	-0.01
14		1,2-Dichlorobenzene	1.550	1.478	4.6	107	-0.01
15		Benzyl Alcohol	1.410	1.310	7.1	95	-0.01
16		2,2'-oxybis(1-Chloropropane	2.802	2.690	4.0	101	-0.01
17		2-Methylphenol	1.225	1.168	4.7	100	0.00
18		Hexachloroethane	0.587	0.553	5.8	102	-0.01
19	P	n-Nitroso-di-n-propylamine	1.304	1.252	4.0	98	-0.01
20		3+4-Methylphenols	1.660	1.623	2.2	104	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
22		Acetophenone	0.598	0.564	5.7	98	-0.01
23	S	Nitrobenzene-d5	0.455	0.412	9.5	98	-0.01
24		Nitrobenzene	0.486	0.498	-2.5	111	-0.01
25		Isophorone	0.881	0.817	7.3	98	-0.01
26	C	2-Nitrophenol	0.198	0.200	-1.0	110	0.00
27		2,4-Dimethylphenol	0.332	0.320	3.6	97	-0.01
28		bis(2-Chloroethoxy)methane	0.502	0.458	8.8	100	-0.01
29	C	2,4-Dichlorophenol	0.320	0.303	5.3	103	0.00
30		1,2,4-Trichlorobenzene	0.348	0.341	2.0	105	-0.01
31		Naphthalene	1.085	1.004	7.5	102	-0.01
32		Benzoic acid	0.229	0.220	3.9	97	0.00
33		4-Chloroaniline	0.468	0.439	6.2	96	-0.01
34	C	Hexachlorobutadiene	0.198	0.186	6.1	104	0.00
35		Caprolactam	0.101	0.101	0.0	103	0.00
36	C	4-Chloro-3-methylphenol	0.359	0.336	6.4	102	-0.01
37		2-Methylnaphthalene	0.725	0.710	2.1	106	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.634	0.585	7.7	100	0.00
40	P	Hexachlorocyclopentadiene	0.301	0.144	52.2#	49#	0.00
41	S	2,4,6-Tribromophenol	0.201	0.176	12.4	94	-0.01
42	C	2,4,6-Trichlorophenol	0.445	0.423	4.9	99	-0.01
43		2,4,5-Trichlorophenol	0.461	0.444	3.7	97	-0.01
44	S	2-Fluorobiphenyl	1.403	1.247	11.1	98	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.754	1.683	4.0	102	-0.01
46		2-Chloronaphthalene	1.336	1.196	10.5	97	-0.01
47		2-Nitroaniline	0.529	0.534	-0.9	100	-0.01
48		Acenaphthylene	2.132	1.906	10.6	97	-0.01
49		Dimethylphthalate	1.625	1.435	11.7	103	-0.01
50		2,6-Dinitrotoluene	0.347	0.299	13.8	90	-0.01
51	C	Acenaphthene	1.254	1.105	11.9	97	-0.01
52		3-Nitroaniline	0.407	0.369	9.3	89	-0.01
53	P	2,4-Dinitrophenol	0.187	0.107	42.8#	55	0.00
54		Dibenzofuran	1.838	1.567	14.7	96	-0.01
55	P	4-Nitrophenol	0.254	0.182	28.3#	74	0.00
56		2,4-Dinitrotoluene	0.440	0.387	12.0	94	-0.01
57		Fluorene	1.451	1.293	10.9	98	-0.01
58		2,3,4,6-Tetrachlorophenol	0.372	0.330	11.3	94	-0.01
59		Diethylphthalate	1.549	1.489	3.9	103	-0.01
60		4-Chlorophenyl-phenylether	0.667	0.631	5.4	100	-0.01
61		4-Nitroaniline	0.407	0.404	0.7	96	-0.01
62		Azobenzene	1.872	1.813	3.2	99	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	98	-0.01
64		4,6-Dinitro-2-methylphenol	0.134	0.111	17.2	75	-0.01
65	c	n-Nitrosodiphenylamine	0.703	0.707	-0.6	98	-0.01
66		4-Bromophenyl-phenylether	0.219	0.212	3.2	96	-0.01
67		Hexachlorobenzene	0.242	0.226	6.6	93	-0.02
68		Atrazine	0.194	0.171	11.9	95	-0.01
69	C	Pentachlorophenol	0.127	0.113	11.0	85	-0.01
70		Phenanthrene	1.184	1.063	10.2	90	-0.02
71		Anthracene	1.192	1.148	3.7	97	-0.01
72		Carbazole	1.071	0.989	7.7	93	-0.01
73		Di-n-butylphthalate	1.269	1.232	2.9	99	-0.02
74	C	Fluoranthene	1.227	1.065	13.2	88	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	74	-0.02
76		Benzidine	0.596	0.547	8.2	69	-0.02
77		Pyrene	1.397	1.502	-7.5	83	-0.02
78	S	Terphenyl-d14	0.839	0.915	-9.1	88	-0.02
79		Butylbenzylphthalate	0.593	0.676	-14.0	90	-0.02
80		Benzo(a)anthracene	1.225	1.191	2.8	75	-0.02
81		3,3'-Dichlorobenzidine	0.383	0.380	0.8	74	-0.02
82		Chrysene	1.184	1.090	7.9	74	-0.02
83		Bis(2-ethylhexyl)phthalate	0.795	0.922	-16.0	92	-0.02
84	c	Di-n-octyl phthalate	1.184	1.442	-21.8#	91	-0.03
85		Indeno(1,2,3-cd)pyrene	0.862	0.949	-10.1	87	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	89	-0.02
87		Benzo(b)fluoranthene	1.283	1.183	7.8	86	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.212	1.090	10.1	84	-0.02
89 C	Benzo(a)pyrene	1.102	1.059	3.9	88	-0.02
90	Dibenzo(a,h)anthracene	0.960	0.909	5.3	89	-0.02
91	Benzo(g,h,i)perylene	0.943	0.783	17.0	79	-0.03

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