Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110715\
Data File : BF082752.D
Acq On : 6 Nov 2015 10:20
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040

Quant Time: Nov 06 17:15:27 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 06 14:39:58 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	121	0.00
2		1,4-Dioxane	0.555	0.523	5.8	112	0.00
3		Pyridine	1.609	1.529	5.0	111	0.00
4		n-Nitrosodimethylamine	0.798	0.783	1.9	121	0.02
5	S	2-Fluorophenol	1.285	1.120	12.8	104	0.00
6		Aniline	2.399	2.352	2.0	125	0.01
7	S	Phenol-d6	1.709	1.627	4.8	119	0.01
8		2-Chlorophenol	1.425	1.301	8.7	111	0.00
9		Benzaldehyde	0.944	0.922	2.3	112	0.00
10	C	Phenol	1.939	1.735	10.5	103	0.01
11		bis(2-Chloroethyl)ether	1.450	1.260	13.1	100	0.00
12		1,3-Dichlorobenzene	1.607	1.337	16.8	99	0.23
13	C	1,4-Dichlorobenzene	1.672	1.646	1.6	131	0.01
14		1,2-Dichlorobenzene	1.521	1.337	12.1	106	0.00
15		Benzyl Alcohol	1.501	1.519	-1.2	127	0.01
16		2,2'-oxybis(1-Chloropropane	2.127	2.030	4.6	119	0.00
17		2-Methylphenol	1.207	1.161	3.8	115	0.01
18		Hexachloroethane	0.598	0.579	3.2	125	0.00
19	P	n-Nitroso-di-n-propylamine	1.256	1.241	1.2	122	0.02
20		3+4-Methylphenols	1.569	1.599	-1.9	137	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	129	0.00
22		Acetophenone	0.573	0.534	6.8	132	0.01
23	S	Nitrobenzene-d5	0.460	0.413	10.2	117	0.01
24		Nitrobenzene	0.470	0.413	12.1	118	0.00
25		Isophorone	0.850	0.783	7.9	124	0.01
26	C	2-Nitrophenol	0.196	0.172	12.2	104	0.00
27		2,4-Dimethylphenol	0.353	0.323	8.5	124	0.01
28		bis(2-Chloroethoxy)methane	0.480	0.404	15.8	106	0.00
29	C	2,4-Dichlorophenol	0.319	0.286	10.3	120	0.00
30		1,2,4-Trichlorobenzene	0.358	0.335	6.4	122	0.00
31		Naphthalene	1.103	0.999	9.4	122	0.01
32		Benzoic acid	0.241	0.238	1.2	125	0.03
33		4-Chloroaniline	0.476	0.430	9.7	118	0.01
34	C	Hexachlorobutadiene	0.224	0.219	2.2	129	0.00
35		Caprolactam	0.100	0.094	6.0	116	0.03
36	C	4-Chloro-3-methylphenol	0.375	0.330	12.0	117	0.00
37		2-Methylnaphthalene	0.714	0.612	14.3	116	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	124	0.01
39		1,2,4,5-Tetrachlorobenzene	0.657	0.621	5.5	108	0.00
40	P	Hexachlorocyclopentadiene	0.353	0.359	-1.7	120	0.00
41	S	2,4,6-Tribromophenol	0.229	0.221	3.5	125	0.00
42	C	2,4,6-Trichlorophenol	0.491	0.489	0.4	128	0.01
43		2,4,5-Trichlorophenol	0.485	0.457	5.8	115	0.01
44	S	2-Fluorobiphenyl	1.395	1.375	1.4	131	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.716	1.521	11.4	102	0.00
46		2-Chloronaphthalene	1.339	1.205	10.0	105	0.00
47		2-Nitroaniline	0.497	0.473	4.8	107	0.01
48		Acenaphthylene	2.098	1.985	5.4	121	0.00
49		Dimethylphthalate	1.639	1.647	-0.5	137	0.01
50		2,6-Dinitrotoluene	0.350	0.369	-5.4	146	0.01
51	C	Acenaphthene	1.330	1.327	0.2	128	0.01
52		3-Nitroaniline	0.385	0.330	14.3	95	0.01
53	P	2,4-Dinitrophenol	0.192	0.180	6.3	100	0.01
54		Dibenzofuran	1.793	1.693	5.6	122	0.00
55	P	4-Nitrophenol	0.295	0.264	10.5	112	0.01
56		2,4-Dinitrotoluene	0.474	0.495	-4.4	143	0.01
57		Fluorene	1.452	1.313	9.6	114	0.00
58		2,3,4,6-Tetrachlorophenol	0.396	0.416	-5.1	140	-0.03
59		Diethylphthalate	1.600	1.641	-2.6	129	0.01
60		4-Chlorophenyl-phenylether	0.726	0.743	-2.3	133	0.01
61		4-Nitroaniline	0.387	0.356	8.0	118	0.01
62		Azobenzene	1.719	1.810	-5.3	131	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	125	0.00
64		4,6-Dinitro-2-methylphenol	0.142	0.155	-9.2	139	0.01
65	c	n-Nitrosodiphenylamine	0.723	0.602	16.7	104	0.00
66		4-Bromophenyl-phenylether	0.241	0.241	0.0	137	0.01
67		Hexachlorobenzene	0.269	0.272	-1.1	139	0.01
68		Atrazine	0.223	0.221	0.9	131	0.01
69	C	Pentachlorophenol	0.163	0.160	1.8	111	0.00
70		Phenanthrene	1.161	1.033	11.0	111	0.00
71		Anthracene	1.223	1.155	5.6	127	0.01
72		Carbazole	1.071	0.888	17.1	102	0.00
73		Di-n-butylphthalate	1.334	1.289	3.4	123	0.00
74	C	Fluoranthene	1.246	1.068	14.3	106	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	117	0.01
76		Benzidine	0.670	0.662	1.2	110	0.00
77		Pyrene	1.373	1.245	9.3	110	0.01
78	S	Terphenyl-d14	0.843	0.745	11.6	102	0.00
79		Butylbenzylphthalate	0.607	0.685	-12.9	141	0.01
80		Benzo(a)anthracene	1.251	1.183	5.4	117	0.00
81		3,3'-Dichlorobenzidine	0.445	0.403	9.4	107	0.01
82		Chrysene	1.146	1.133	1.1	126	0.01
83		Bis(2-ethylhexyl)phthalate	0.831	0.925	-11.3	138	0.01
84	c	Di-n-octyl phthalate	1.385	1.397	-0.9	120	0.01
85		Indeno(1,2,3-cd)pyrene	0.987	0.929	5.9	117	0.03
86	I	Perylene-d12	1.000	1.000	0.0	122	0.02
87		Benzo(b)fluoranthene	1.284	1.172	8.7	123	0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.139	1.172	-2.9	120	-0.01
89 C	Benzo(a)pyrene	1.112	1.044	6.1	117	0.02
90	Dibenzo(a,h)anthracene	0.942	0.885	6.1	117	0.03
91	Benzo(g,h,i)perylene	0.902	0.823	8.8	116	0.03

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