Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020116\
Data File : BF084708.D
Acq On : 1 Feb 2016 14:37
Operator : SJ/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF020116

Quant Time: Feb 01 15:07:18 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 01 14:55:11 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	102	-0.01
2		1,4-Dioxane	0.562	0.511	9.1	95	-0.01
3		Pyridine	1.465	1.329	9.3	98	-0.02
4		n-Nitrosodimethylamine	0.758	0.726	4.2	97	-0.02
5	S	2-Fluorophenol	1.284	1.176	8.4	100	-0.01
6		Aniline	1.948	1.896	2.7	101	-0.01
7	S	Phenol-d6	1.592	1.428	10.3	98	-0.01
8		2-Chlorophenol	1.453	1.419	2.3	99	-0.01
9		Benzaldehyde	0.940	0.955	-1.6	102	-0.01
10	C	Phenol	1.783	1.511	15.3	99	-0.01
11		bis(2-Chloroethyl)ether	1.391	1.240	10.9	99	-0.01
12		1,3-Dichlorobenzene	1.536	1.404	8.6	99	-0.01
13	C	1,4-Dichlorobenzene	1.535	1.448	5.7	101	-0.01
14		1,2-Dichlorobenzene	1.430	1.401	2.0	98	-0.01
15		Benzyl Alcohol	1.099	1.044	5.0	100	-0.01
16		2,2'-oxybis(1-Chloropropane	2.339	2.200	5.9	102	-0.01
17		2-Methylphenol	1.149	1.166	-1.5	99	-0.01
18		Hexachloroethane	0.591	0.580	1.9	99	-0.01
19	P	n-Nitroso-di-n-propylamine	0.998	0.896	10.2	98	-0.01
20		3+4-Methylphenols	1.427	1.368	4.1	100	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	101	-0.02
22		Acetophenone	0.527	0.510	3.2	100	-0.01
23	S	Nitrobenzene-d5	0.355	0.360	-1.4	101	-0.01
24		Nitrobenzene	0.380	0.348	8.4	102	-0.01
25		Isophorone	0.690	0.691	-0.1	100	-0.01
26	C	2-Nitrophenol	0.188	0.178	5.3	97	-0.01
27		2,4-Dimethylphenol	0.326	0.299	8.3	99	-0.01
28		bis(2-Chloroethoxy)methane	0.410	0.395	3.7	101	-0.01
29	C	2,4-Dichlorophenol	0.279	0.285	-2.2	99	-0.01
30		1,2,4-Trichlorobenzene	0.315	0.307	2.5	101	-0.01
31		Naphthalene	1.003	0.924	7.9	99	-0.01
32		Benzoic acid	0.209	0.224	-7.2	108	-0.01
33		4-Chloroaniline	0.415	0.379	8.7	101	-0.01
34	C	Hexachlorobutadiene	0.182	0.176	3.3	96	-0.01
35		Caprolactam	0.086	0.096	-11.6	99	0.00
36	C	4-Chloro-3-methylphenol	0.318	0.302	5.0	103	-0.01
37		2-Methylnaphthalene	0.628	0.659	-4.9	101	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	100	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.635	0.589	7.2	102	-0.02
40	P	Hexachlorocyclopentadiene	0.223	0.240	-7.6	107	-0.01
41	S	2,4,6-Tribromophenol	0.209	0.217	-3.8	99	-0.01
42	C	2,4,6-Trichlorophenol	0.409	0.404	1.2	97	-0.01
43		2,4,5-Trichlorophenol	0.416	0.386	7.2	97	-0.01
44	S	2-Fluorobiphenyl	1.321	1.202	9.0	102	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.786	1.821	-2.0	102	-0.01
46		2-Chloronaphthalene	1.316	1.268	3.6	99	-0.01
47		2-Nitroaniline	0.417	0.376	9.8	103	-0.01
48		Acenaphthylene	1.996	2.005	-0.5	101	-0.01
49		Dimethylphthalate	1.523	1.538	-1.0	100	-0.01
50		2,6-Dinitrotoluene	0.333	0.345	-3.6	99	-0.01
51	C	Acenaphthene	1.299	1.271	2.2	99	-0.01
52		3-Nitroaniline	0.374	0.384	-2.7	107	0.00
53	P	2,4-Dinitrophenol	0.133	0.152	-14.3	107	0.00
54		Dibenzofuran	1.587	1.608	-1.3	101	-0.01
55	P	4-Nitrophenol	0.275	0.294	-6.9	98	0.00
56		2,4-Dinitrotoluene	0.424	0.415	2.1	97	-0.01
57		Fluorene	1.326	1.376	-3.8	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.317	0.299	5.7	106	0.00
59		Diethylphthalate	1.487	1.499	-0.8	105	0.00
60		4-Chlorophenyl-phenylether	0.621	0.568	8.5	98	-0.01
61		4-Nitroaniline	0.364	0.346	4.9	98	-0.01
62		Azobenzene	1.430	1.394	2.5	99	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	-0.01
64		4,6-Dinitro-2-methylphenol	0.123	0.137	-11.4	105	0.00
65	c	n-Nitrosodiphenylamine	0.635	0.670	-5.5	98	-0.01
66		4-Bromophenyl-phenylether	0.221	0.235	-6.3	99	-0.01
67		Hexachlorobenzene	0.241	0.235	2.5	102	-0.01
68		Atrazine	0.201	0.200	0.5	106	-0.01
69	C	Pentachlorophenol	0.113	0.121	-7.1	103	0.00
70		Phenanthrene	1.109	1.047	5.6	101	-0.01
71		Anthracene	1.118	1.179	-5.5	99	-0.01
72		Carbazole	0.975	1.062	-8.9	101	-0.01
73		Di-n-butylphthalate	1.241	1.192	3.9	100	-0.01
74	C	Fluoranthene	1.123	1.161	-3.4	99	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	99	0.00
76		Benzidine	0.597	0.649	-8.7	90	-0.01
77		Pyrene	1.531	1.615	-5.5	100	-0.01
78	S	Terphenyl-d14	0.883	0.843	4.5	99	-0.01
79		Butylbenzylphthalate	0.695	0.734	-5.6	98	-0.01
80		Benzo(a)anthracene	1.241	1.249	-0.6	99	0.00
81		3,3'-Dichlorobenzidine	0.440	0.485	-10.2	99	-0.01
82		Chrysene	1.204	1.272	-5.6	101	-0.01
83		Bis(2-ethylhexyl)phthalate	0.864	0.871	-0.8	97	-0.01
84	c	Di-n-octyl phthalate	1.426	1.476	-3.5	99	0.00
85		Indeno(1,2,3-cd)pyrene	0.947	0.911	3.8	97	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	99	-0.01
87		Benzo(b)fluoranthene	1.344	1.345	-0.1	98	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.111	1.110	0.1	102	-0.01
89 C	Benzo(a)pyrene	1.145	1.149	-0.3	99	-0.01
90	Dibenzo(a,h)anthracene	0.964	0.933	3.2	97	-0.01
91	Benzo(g,h,i)perylene	0.971	0.922	5.0	96	-0.01

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