Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
Data File : BF085713.D
Acq On : 24 Mar 2016 13:39
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF032416

Quant Time: Mar 25 14:41:11 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 24 14:15:39 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	98	0.00
2		1,4-Dioxane	0.573	0.584	-1.9	101	0.00
3		Pyridine	1.375	1.558	-13.3	105	0.00
4		n-Nitrosodimethylamine	0.550	0.535	2.7	91	0.01
5	S	2-Fluorophenol	1.201	1.191	0.8	102	0.00
6		Aniline	2.055	2.029	1.3	101	0.00
7	S	Phenol-d6	1.527	1.473	3.5	100	0.00
8		2-Chlorophenol	1.395	1.442	-3.4	101	0.00
9		Benzaldehyde	0.920	0.906	1.5	101	0.00
10	C	Phenol	1.730	1.723	0.4	102	0.00
11		bis(2-Chloroethyl)ether	1.331	1.322	0.7	101	0.00
12		1,3-Dichlorobenzene	1.503	1.546	-2.9	101	0.00
13	C	1,4-Dichlorobenzene	1.524	1.505	1.2	103	0.00
14		1,2-Dichlorobenzene	1.420	1.455	-2.5	101	0.00
15		Benzyl Alcohol	1.019	1.078	-5.8	100	0.00
16		2,2'-oxybis(1-Chloropropane	1.766	1.795	-1.6	102	0.00
17		2-Methylphenol	1.117	1.169	-4.7	101	0.00
18		Hexachloroethane	0.558	0.580	-3.9	105	0.00
19	P	n-Nitroso-di-n-propylamine	0.914	0.973	-6.5	100	0.00
20		3+4-Methylphenols	1.343	1.344	-0.1	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
22		Acetophenone	0.466	0.445	4.5	101	0.00
23	S	Nitrobenzene-d5	0.355	0.354	0.3	100	0.00
24		Nitrobenzene	0.366	0.344	6.0	101	0.00
25		Isophorone	0.684	0.675	1.3	104	0.00
26	C	2-Nitrophenol	0.183	0.180	1.6	102	0.00
27		2,4-Dimethylphenol	0.320	0.343	-7.2	102	0.00
28		bis(2-Chloroethoxy)methane	0.390	0.402	-3.1	101	0.00
29	C	2,4-Dichlorophenol	0.269	0.258	4.1	99	0.00
30		1,2,4-Trichlorobenzene	0.299	0.301	-0.7	104	0.00
31		Naphthalene	1.008	0.965	4.3	102	0.00
32		Benzoic acid	0.213	0.225	-5.6	96	0.00
33		4-Chloroaniline	0.411	0.403	1.9	100	0.00
34	C	Hexachlorobutadiene	0.162	0.162	0.0	103	0.00
35		Caprolactam	0.095	0.087	8.4	99	0.00
36	C	4-Chloro-3-methylphenol	0.314	0.292	7.0	101	0.00
37		2-Methylnaphthalene	0.635	0.610	3.9	100	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
39		1,2,4,5-Tetrachlorobenzene	0.593	0.627	-5.7	102	0.00
40	P	Hexachlorocyclopentadiene	0.196	0.209	-6.6	103	-0.01
41	S	2,4,6-Tribromophenol	0.190	0.198	-4.2	100	0.00
42	C	2,4,6-Trichlorophenol	0.401	0.405	-1.0	101	0.00
43		2,4,5-Trichlorophenol	0.420	0.439	-4.5	103	0.00
44	S	2-Fluorobiphenyl	1.197	1.228	-2.6	98	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.688	1.804	-6.9	104	0.00
46		2-Chloronaphthalene	1.241	1.338	-7.8	101	0.00
47		2-Nitroaniline	0.379	0.355	6.3	98	0.00
48		Acenaphthylene	2.022	2.122	-4.9	100	0.00
49		Dimethylphthalate	1.517	1.430	5.7	99	0.00
50		2,6-Dinitrotoluene	0.327	0.359	-9.8	103	0.00
51	C	Acenaphthene	1.235	1.284	-4.0	108	0.00
52		3-Nitroaniline	0.374	0.363	2.9	102	0.00
53	P	2,4-Dinitrophenol	0.140	0.164	-17.1	104	0.00
54		Dibenzofuran	1.599	1.638	-2.4	101	0.00
55	P	4-Nitrophenol	0.244	0.258	-5.7	99	0.00
56		2,4-Dinitrotoluene	0.415	0.445	-7.2	96	0.00
57		Fluorene	1.329	1.415	-6.5	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.292	0.290	0.7	99	0.00
59		Diethylphthalate	1.499	1.482	1.1	100	0.00
60		4-Chlorophenyl-phenylether	0.650	0.617	5.1	102	0.00
61		4-Nitroaniline	0.358	0.387	-8.1	100	0.00
62		Azobenzene	1.441	1.497	-3.9	104	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
64		4,6-Dinitro-2-methylphenol	0.117	0.131	-12.0	102	0.00
65	c	n-Nitrosodiphenylamine	0.633	0.656	-3.6	101	0.00
66		4-Bromophenyl-phenylether	0.205	0.208	-1.5	100	0.00
67		Hexachlorobenzene	0.214	0.224	-4.7	99	0.00
68		Atrazine	0.214	0.211	1.4	100	0.00
69	C	Pentachlorophenol	0.105	0.116	-10.5	100	0.00
70		Phenanthrene	1.027	1.115	-8.6	100	0.00
71		Anthracene	1.078	1.068	0.9	103	0.00
72		Carbazole	0.905	0.936	-3.4	98	0.00
73		Di-n-butylphthalate	1.242	1.248	-0.5	102	0.00
74	C	Fluoranthene	1.092	1.073	1.7	100	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	93	0.00
76		Benzidine	0.704	0.720	-2.3	94	0.00
77		Pyrene	1.644	1.569	4.6	99	0.00
78	S	Terphenyl-d14	0.972	0.947	2.6	102	0.00
79		Butylbenzylphthalate	0.688	0.702	-2.0	98	0.00
80		Benzo(a)anthracene	1.140	1.125	1.3	96	0.00
81		3,3'-Dichlorobenzidine	0.791	0.722	8.7	94	0.00
82		Chrysene	1.149	1.102	4.1	95	-0.01
83		Bis(2-ethylhexyl)phthalate	0.824	0.832	-1.0	95	0.00
84	c	Di-n-octyl phthalate	1.226	1.226	0.0	92	0.00
85		Indeno(1,2,3-cd)pyrene	0.887	0.905	-2.0	97	0.00
86	I	Perylene-d12	1.000	1.000	0.0	92	0.00
87		Benzo(b)fluoranthene	1.260	1.360	-7.9	98	0.00

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88	Benzo(k)fluoranthene	1.136	1.057	7.0	93	0.00
89 C	Benzo(a)pyrene	1.119	1.141	-2.0	96	0.00
90	Dibenzo(a,h)anthracene	1.039	1.069	-2.9	97	0.00
91	Benzo(g,h,i)perylene	1.052	1.079	-2.6	97	-0.01

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