Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020316\
Data File : BF084804.D
Acq On : 4 Feb 2016 00:24
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 04 04:41:39 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 03 14:16:01 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	87	0.00
2		1,4-Dioxane	0.562	0.576	-2.5	92	-0.01
3		Pyridine	1.465	1.591	-8.6	100	-0.01
4		n-Nitrosodimethylamine	0.758	0.725	4.4	83	-0.01
5	S	2-Fluorophenol	1.284	1.361	-6.0	99	0.00
6		Aniline	1.948	2.053	-5.4	94	0.00
7	S	Phenol-d6	1.592	1.572	1.3	93	0.00
8		2-Chlorophenol	1.453	1.537	-5.8	91	0.00
9		Benzaldehyde	0.940	0.949	-1.0	87	-0.01
10	C	Phenol	1.783	1.724	3.3	97	0.00
11		bis(2-Chloroethyl)ether	1.391	1.419	-2.0	97	0.00
12		1,3-Dichlorobenzene	1.536	1.593	-3.7	96	0.00
13	C	1,4-Dichlorobenzene	1.535	1.639	-6.8	98	0.00
14		1,2-Dichlorobenzene	1.430	1.349	5.7	81	0.00
15		Benzyl Alcohol	1.099	1.134	-3.2	93	0.00
16		2,2'-oxybis(1-Chloropropane	2.339	2.310	1.2	92	0.00
17		2-Methylphenol	1.149	1.187	-3.3	86	0.00
18		Hexachloroethane	0.591	0.588	0.5	86	0.00
19	P	n-Nitroso-di-n-propylamine	0.998	0.989	0.9	93	0.00
20		3+4-Methylphenols	1.427	1.401	1.8	88	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	90	0.00
22		Acetophenone	0.527	0.506	4.0	89	0.00
23	S	Nitrobenzene-d5	0.355	0.372	-4.8	93	0.00
24		Nitrobenzene	0.380	0.340	10.5	90	0.00
25		Isophorone	0.690	0.681	1.3	88	0.00
26	C	2-Nitrophenol	0.188	0.194	-3.2	95	0.00
27		2,4-Dimethylphenol	0.326	0.292	10.4	87	-0.01
28		bis(2-Chloroethoxy)methane	0.410	0.376	8.3	86	-0.01
29	C	2,4-Dichlorophenol	0.279	0.276	1.1	86	0.00
30		1,2,4-Trichlorobenzene	0.315	0.299	5.1	88	0.00
31		Naphthalene	1.003	0.944	5.9	91	0.00
32		Benzoic acid	0.209	0.231	-10.5	100	0.00
33		4-Chloroaniline	0.415	0.398	4.1	95	0.00
34	C	Hexachlorobutadiene	0.182	0.181	0.5	89	0.00
35		Caprolactam	0.086	0.094	-9.3	87	0.00
36	C	4-Chloro-3-methylphenol	0.318	0.300	5.7	91	0.00
37		2-Methylnaphthalene	0.628	0.645	-2.7	89	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	84	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.635	0.591	6.9	87	0.00
40	P	Hexachlorocyclopentadiene	0.223	0.262	-17.5	99	0.00
41	S	2,4,6-Tribromophenol	0.209	0.209	0.0	81	0.00
42	C	2,4,6-Trichlorophenol	0.409	0.400	2.2	81	-0.01
43		2,4,5-Trichlorophenol	0.416	0.416	0.0	88	0.00
44	S	2-Fluorobiphenyl	1.321	1.364	-3.3	98	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.786	1.704	4.6	80	-0.01
46		2-Chloronaphthalene	1.316	1.226	6.8	81	0.00
47		2-Nitroaniline	0.417	0.375	10.1	87	0.00
48		Acenaphthylene	1.996	1.838	7.9	78	-0.01
49		Dimethylphthalate	1.523	1.539	-1.1	85	0.00
50		2,6-Dinitrotoluene	0.333	0.369	-10.8	89	0.00
51	C	Acenaphthene	1.299	1.223	5.9	81	-0.01
52		3-Nitroaniline	0.374	0.339	9.4	80	-0.01
53	P	2,4-Dinitrophenol	0.133	0.158	-18.8	94	0.00
54		Dibenzofuran	1.587	1.529	3.7	81	-0.01
55	P	4-Nitrophenol	0.275	0.272	1.1	76	0.00
56		2,4-Dinitrotoluene	0.424	0.466	-9.9	92	0.00
57		Fluorene	1.326	1.236	6.8	77	0.00
58		2,3,4,6-Tetrachlorophenol	0.317	0.333	-5.0	100	0.00
59		Diethylphthalate	1.487	1.435	3.5	85	-0.01
60		4-Chlorophenyl-phenylether	0.621	0.642	-3.4	94	0.00
61		4-Nitroaniline	0.364	0.376	-3.3	90	0.00
62		Azobenzene	1.430	1.548	-8.3	93	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
64		4,6-Dinitro-2-methylphenol	0.123	0.130	-5.7	90	0.00
65	c	n-Nitrosodiphenylamine	0.635	0.664	-4.6	88	0.00
66		4-Bromophenyl-phenylether	0.221	0.227	-2.7	87	0.00
67		Hexachlorobenzene	0.241	0.234	2.9	92	0.00
68		Atrazine	0.201	0.184	8.5	89	0.00
69	C	Pentachlorophenol	0.113	0.113	0.0	87	0.00
70		Phenanthrene	1.109	1.056	4.8	92	0.00
71		Anthracene	1.118	1.116	0.2	85	0.00
72		Carbazole	0.975	0.906	7.1	78	-0.01
73		Di-n-butylphthalate	1.241	1.157	6.8	88	0.00
74	C	Fluoranthene	1.123	1.099	2.1	85	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	90	0.00
76		Benzidine	0.597	0.708	-18.6	89	0.00
77		Pyrene	1.531	1.395	8.9	79	0.00
78	S	Terphenyl-d14	0.883	0.907	-2.7	97	0.00
79		Butylbenzylphthalate	0.695	0.652	6.2	79	-0.01
80		Benzo(a)anthracene	1.241	1.184	4.6	86	0.00
81		3,3'-Dichlorobenzidine	0.440	0.392	10.9	73	-0.01
82		Chrysene	1.204	1.028	14.6	75	0.00
83		Bis(2-ethylhexyl)phthalate	0.864	0.871	-0.8	88	0.00
84	c	Di-n-octyl phthalate	1.426	1.322	7.3	81	0.00
85		Indeno(1,2,3-cd)pyrene	0.947	0.905	4.4	88	0.00
86	I	Perylene-d12	1.000	1.000	0.0	79	0.00
87		Benzo(b)fluoranthene	1.344	1.252	6.8	73	0.00

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88	Benzo(k)fluoranthene	1.111	1.179	-6.1	86	0.00
89 C	Benzo(a)pyrene	1.145	1.117	2.4	76	0.00
90	Dibenzo(a,h)anthracene	0.964	1.054	-9.3	88	0.00
91	Benzo(g,h,i)perylene	0.971	1.092	-12.5	90	0.00

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