Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011516\
Data File : BF084486.D
Acq On : 15 Jan 2016 11:08
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 15 11:34:37 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 14 14:15:52 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	84	-0.01
2		1,4-Dioxane	0.586	0.611	-4.3	83	-0.01
3		Pyridine	1.633	1.722	-5.5	81	-0.02
4		n-Nitrosodimethylamine	0.838	0.765	8.7	72	-0.01
5	S	2-Fluorophenol	1.236	1.263	-2.2	91	-0.01
6		Aniline	2.272	2.053	9.6	73	-0.01
7	S	Phenol-d6	1.553	1.563	-0.6	83	-0.01
8		2-Chlorophenol	1.403	1.472	-4.9	89	-0.01
9		Benzaldehyde	1.011	0.914	9.6	76	-0.01
10	C	Phenol	1.766	1.719	2.7	75	-0.01
11		bis(2-Chloroethyl)ether	1.368	1.504	-9.9	96	-0.01
12		1,3-Dichlorobenzene	1.447	1.467	-1.4	87	-0.01
13	C	1,4-Dichlorobenzene	1.451	1.511	-4.1	83	-0.01
14		1,2-Dichlorobenzene	1.390	1.272	8.5	82	-0.01
15		Benzyl Alcohol	1.306	1.139	12.8	69	-0.01
16		2,2'-oxybis(1-Chloropropane	2.491	2.254	9.5	78	-0.01
17		2-Methylphenol	1.176	1.224	-4.1	81	-0.01
18		Hexachloroethane	0.580	0.560	3.4	78	-0.01
19	P	n-Nitroso-di-n-propylamine	1.051	0.940	10.6	77	-0.01
20		3+4-Methylphenols	1.382	1.324	4.2	85	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	84	-0.01
22		Acetophenone	0.481	0.473	1.7	77	-0.01
23	S	Nitrobenzene-d5	0.359	0.325	9.5	77	-0.01
24		Nitrobenzene	0.364	0.387	-6.3	80	-0.01
25		Isophorone	0.687	0.638	7.1	77	-0.01
26	C	2-Nitrophenol	0.166	0.187	-12.7	96	-0.01
27		2,4-Dimethylphenol	0.336	0.322	4.2	75	0.00
28		bis(2-Chloroethoxy)methane	0.420	0.430	-2.4	92	-0.01
29	C	2,4-Dichlorophenol	0.280	0.276	1.4	84	-0.01
30		1,2,4-Trichlorobenzene	0.289	0.294	-1.7	82	-0.01
31		Naphthalene	0.907	0.885	2.4	82	-0.01
32		Benzoic acid	0.198	0.174	12.1	70	0.00
33		4-Chloroaniline	0.416	0.429	-3.1	89	-0.01
34	C	Hexachlorobutadiene	0.166	0.154	7.2	77	-0.01
35		Caprolactam	0.094	0.087	7.4	68	0.00
36	C	4-Chloro-3-methylphenol	0.313	0.302	3.5	84	0.00
37		2-Methylnaphthalene	0.651	0.631	3.1	84	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	83	0.00
39		1,2,4,5-Tetrachlorobenzene	0.575	0.540	6.1	78	-0.01
40	P	Hexachlorocyclopentadiene	0.347	0.286	17.6	67	-0.01
41	S	2,4,6-Tribromophenol	0.202	0.195	3.5	75	-0.01
42	C	2,4,6-Trichlorophenol	0.430	0.377	12.3	75	-0.01
43		2,4,5-Trichlorophenol	0.412	0.405	1.7	79	-0.01
44	S	2-Fluorobiphenyl	1.317	1.188	9.8	85	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.590	1.496	5.9	84	-0.01
46		2-Chloronaphthalene	1.249	1.152	7.8	85	-0.01
47		2-Nitroaniline	0.412	0.408	1.0	85	-0.01
48		Acenaphthylene	1.845	1.876	-1.7	80	0.00
49		Dimethylphthalate	1.430	1.243	13.1	69	-0.01
50		2,6-Dinitrotoluene	0.314	0.293	6.7	70	-0.01
51	C	Acenaphthene	1.237	1.270	-2.7	79	0.00
52		3-Nitroaniline	0.389	0.339	12.9	67	-0.01
53	P	2,4-Dinitrophenol	0.093	0.151	-62.4#	94	-0.01
54		Dibenzofuran	1.812	1.636	9.7	78	0.00
55	P	4-Nitrophenol	0.282	0.260	7.8	64	-0.01
56		2,4-Dinitrotoluene	0.397	0.418	-5.3	75	0.00
57		Fluorene	1.400	1.268	9.4	80	0.00
58		2,3,4,6-Tetrachlorophenol	0.352	0.314	10.8	70	-0.01
59		Diethylphthalate	1.410	1.247	11.6	72	-0.01
60		4-Chlorophenyl-phenylether	0.564	0.535	5.1	79	-0.01
61		4-Nitroaniline	0.346	0.363	-4.9	89	0.00
62		Azobenzene	1.546	1.425	7.8	75	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	82	-0.01
64		4,6-Dinitro-2-methylphenol	0.109	0.113	-3.7	82	-0.01
65	c	n-Nitrosodiphenylamine	0.639	0.659	-3.1	84	0.00
66		4-Bromophenyl-phenylether	0.215	0.215	0.0	77	-0.01
67		Hexachlorobenzene	0.242	0.233	3.7	83	0.00
68		Atrazine	0.209	0.187	10.5	65	-0.01
69	C	Pentachlorophenol	0.126	0.117	7.1	67	0.00
70		Phenanthrene	1.046	1.094	-4.6	79	0.00
71		Anthracene	1.026	0.953	7.1	79	-0.01
72		Carbazole	1.003	0.918	8.5	72	-0.01
73		Di-n-butylphthalate	1.111	1.039	6.5	75	-0.01
74	C	Fluoranthene	1.093	1.041	4.8	81	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	76	-0.01
76		Benzidine	0.717	0.634	11.6	66	-0.01
77		Pyrene	1.569	1.503	4.2	82	-0.01
78	S	Terphenyl-d14	0.896	0.805	10.2	79	-0.01
79		Butylbenzylphthalate	0.679	0.648	4.6	71	-0.01
80		Benzo(a)anthracene	1.174	1.165	0.8	80	-0.01
81		3,3'-Dichlorobenzidine	0.410	0.478	-16.6	88	0.00
82		Chrysene	1.128	1.198	-6.2	82	0.00
83		Bis(2-ethylhexyl)phthalate	0.858	0.783	8.7	72	-0.01
84	c	Di-n-octyl phthalate	1.449	1.425	1.7	71	0.00
85		Indeno(1,2,3-cd)pyrene	0.895	1.027	-14.7	89	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	88	-0.01
87		Benzo(b)fluoranthene	1.308	1.201	8.2	77	-0.01

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88	Benzo(k)fluoranthene	1.070	1.100	-2.8	91	-0.01
89 C	Benzo(a)pyrene	1.110	1.086	2.2	82	-0.01
90	Dibenzo(a,h)anthracene	0.955	0.925	3.1	89	0.00
91	Benzo(g,h,i)perylene	0.989	0.964	2.5	92	-0.01

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