Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF123115\
Data File : BF084206.D
Acq On : 31 Dec 2015 20:31
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 02 04:18:30 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 31 14:15:57 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	90	0.00
2		1,4-Dioxane	0.586	0.622	-6.1	91	-0.01
3		Pyridine	1.633	1.698	-4.0	86	-0.02
4		n-Nitrosodimethylamine	0.838	0.895	-6.8	91	-0.02
5	S	2-Fluorophenol	1.236	1.226	0.8	95	0.00
6		Aniline	2.272	2.398	-5.5	92	0.00
7	S	Phenol-d6	1.553	1.667	-7.3	96	0.00
8		2-Chlorophenol	1.403	1.409	-0.4	92	0.00
9		Benzaldehyde	1.011	0.984	2.7	88	-0.01
10	C	Phenol	1.766	2.140	-21.2#	101	0.00
11		bis(2-Chloroethyl)ether	1.368	1.298	5.1	89	0.00
12		1,3-Dichlorobenzene	1.447	1.479	-2.2	95	0.00
13	C	1,4-Dichlorobenzene	1.451	1.571	-8.3	93	0.00
14		1,2-Dichlorobenzene	1.390	1.330	4.3	92	-0.01
15		Benzyl Alcohol	1.306	1.418	-8.6	93	0.00
16		2,2'-oxybis(1-Chloropropane	2.491	2.484	0.3	92	0.00
17		2-Methylphenol	1.176	1.259	-7.1	90	-0.01
18		Hexachloroethane	0.580	0.608	-4.8	91	0.00
19	P	n-Nitroso-di-n-propylamine	1.051	0.991	5.7	87	0.00
20		3+4-Methylphenols	1.382	1.345	2.7	93	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
22		Acetophenone	0.481	0.501	-4.2	90	0.00
23	S	Nitrobenzene-d5	0.359	0.339	5.6	89	0.00
24		Nitrobenzene	0.364	0.385	-5.8	88	0.00
25		Isophorone	0.687	0.685	0.3	92	0.00
26	C	2-Nitrophenol	0.166	0.165	0.6	94	0.00
27		2,4-Dimethylphenol	0.336	0.343	-2.1	89	0.00
28		bis(2-Chloroethoxy)methane	0.420	0.377	10.2	90	0.00
29	C	2,4-Dichlorophenol	0.280	0.266	5.0	90	0.00
30		1,2,4-Trichlorobenzene	0.289	0.305	-5.5	94	0.00
31		Naphthalene	0.907	0.907	0.0	93	0.00
32		Benzoic acid	0.198	0.193	2.5	86	0.00
33		4-Chloroaniline	0.416	0.407	2.2	93	0.00
34	C	Hexachlorobutadiene	0.166	0.168	-1.2	94	0.00
35		Caprolactam	0.094	0.086	8.5	75	-0.01
36	C	4-Chloro-3-methylphenol	0.313	0.311	0.6	96	0.00
37		2-Methylnaphthalene	0.651	0.599	8.0	89	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
39		1,2,4,5-Tetrachlorobenzene	0.575	0.604	-5.0	93	0.00
40	P	Hexachlorocyclopentadiene	0.347	0.374	-7.8	94	0.00
41	S	2,4,6-Tribromophenol	0.202	0.205	-1.5	85	0.00
42	C	2,4,6-Trichlorophenol	0.430	0.399	7.2	84	0.00
43		2,4,5-Trichlorophenol	0.412	0.448	-8.7	93	0.00
44	S	2-Fluorobiphenyl	1.317	1.150	12.7	88	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.590	1.494	6.0	90	0.00
46		2-Chloronaphthalene	1.249	1.148	8.1	91	0.00
47		2-Nitroaniline	0.412	0.387	6.1	87	0.00
48		Acenaphthylene	1.845	1.914	-3.7	88	0.00
49		Dimethylphthalate	1.430	1.527	-6.8	91	0.00
50		2,6-Dinitrotoluene	0.314	0.354	-12.7	91	0.00
51	C	Acenaphthene	1.237	1.334	-7.8	89	0.00
52		3-Nitroaniline	0.389	0.348	10.5	73	-0.01
53	P	2,4-Dinitrophenol	0.093	0.139	-49.5#	93	0.00
54		Dibenzofuran	1.812	1.714	5.4	87	0.00
55	P	4-Nitrophenol	0.282	0.306	-8.5	80	0.00
56		2,4-Dinitrotoluene	0.397	0.472	-18.9	90	0.00
57		Fluorene	1.400	1.216	13.1	82	0.00
58		2,3,4,6-Tetrachlorophenol	0.352	0.361	-2.6	86	0.00
59		Diethylphthalate	1.410	1.465	-3.9	90	0.00
60		4-Chlorophenyl-phenylether	0.564	0.608	-7.8	96	0.00
61		4-Nitroaniline	0.346	0.352	-1.7	92	0.00
62		Azobenzene	1.546	1.629	-5.4	92	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	88	0.00
64		4,6-Dinitro-2-methylphenol	0.109	0.105	3.7	81	0.00
65	c	n-Nitrosodiphenylamine	0.639	0.594	7.0	81	0.00
66		4-Bromophenyl-phenylether	0.215	0.235	-9.3	90	0.00
67		Hexachlorobenzene	0.242	0.228	5.8	87	0.00
68		Atrazine	0.209	0.204	2.4	76	-0.01
69	C	Pentachlorophenol	0.126	0.140	-11.1	86	0.00
70		Phenanthrene	1.046	1.071	-2.4	83	-0.01
71		Anthracene	1.026	1.042	-1.6	93	0.00
72		Carbazole	1.003	0.906	9.7	76	-0.01
73		Di-n-butylphthalate	1.111	1.196	-7.7	92	0.00
74	C	Fluoranthene	1.093	0.971	11.2	81	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	81	0.00
76		Benzidine	0.717	0.696	2.9	77	0.00
77		Pyrene	1.569	1.415	9.8	82	0.00
78	S	Terphenyl-d14	0.896	0.772	13.8	81	0.00
79		Butylbenzylphthalate	0.679	0.694	-2.2	81	0.00
80		Benzo(a)anthracene	1.174	1.109	5.5	81	0.00
81		3,3'-Dichlorobenzidine	0.410	0.418	-2.0	81	0.00
82		Chrysene	1.128	0.891	21.0	65	-0.01
83		Bis(2-ethylhexyl)phthalate	0.858	0.811	5.5	79	0.00
84	c	Di-n-octyl phthalate	1.449	1.270	12.4	67	-0.01
85		Indeno(1,2,3-cd)pyrene	0.895	0.979	-9.4	90	0.00
86	I	Perylene-d12	1.000	1.000	0.0	77	0.00
87		Benzo(b)fluoranthene	1.308	1.425	-8.9	80	-0.01

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88	Benzo(k)fluoranthene	1.070	0.918	14.2	66	0.00
89 C	Benzo(a)pyrene	1.110	1.133	-2.1	75	-0.01
90	Dibenzo(a,h)anthracene	0.955	1.071	-12.1	90	0.00
91	Benzo(g,h,i)perylene	0.989	1.109	-12.1	92	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1