Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011516\
Data File : BF084509.D
Acq On : 15 Jan 2016 23:45
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 16 03:19:19 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jan 16 01:26:38 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	-0.01
2		1,4-Dioxane	0.586	0.608	-3.8	86	0.00
3		Pyridine	1.633	1.780	-9.0	87	-0.02
4		n-Nitrosodimethylamine	0.838	0.754	10.0	73	-0.01
5	S	2-Fluorophenol	1.236	1.190	3.7	89	-0.01
6		Aniline	2.272	1.957	13.9	72	-0.01
7	S	Phenol-d6	1.553	1.577	-1.5	87	0.00
8		2-Chlorophenol	1.403	1.477	-5.3	92	-0.01
9		Benzaldehyde	1.011	0.932	7.8	80	-0.01
10	C	Phenol	1.766	1.837	-4.0	83	-0.01
11		bis(2-Chloroethyl)ether	1.368	1.482	-8.3	98	-0.01
12		1,3-Dichlorobenzene	1.447	1.456	-0.6	89	-0.01
13	C	1,4-Dichlorobenzene	1.451	1.490	-2.7	85	-0.01
14		1,2-Dichlorobenzene	1.390	1.298	6.6	86	-0.01
15		Benzyl Alcohol	1.306	1.161	11.1	73	-0.01
16		2,2'-oxybis(1-Chloropropane	2.491	2.200	11.7	78	-0.01
17		2-Methylphenol	1.176	1.155	1.8	79	-0.01
18		Hexachloroethane	0.580	0.556	4.1	80	-0.01
19	P	n-Nitroso-di-n-propylamine	1.051	0.957	8.9	81	-0.01
20		3+4-Methylphenols	1.382	1.340	3.0	89	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	85	-0.01
22		Acetophenone	0.481	0.464	3.5	77	-0.01
23	S	Nitrobenzene-d5	0.359	0.333	7.2	81	-0.01
24		Nitrobenzene	0.364	0.372	-2.2	78	-0.01
25		Isophorone	0.687	0.641	6.7	79	-0.01
26	C	2-Nitrophenol	0.166	0.191	-15.1	100	-0.01
27		2,4-Dimethylphenol	0.336	0.335	0.3	80	0.00
28		bis(2-Chloroethoxy)methane	0.420	0.438	-4.3	96	-0.01
29	C	2,4-Dichlorophenol	0.280	0.266	5.0	83	-0.01
30		1,2,4-Trichlorobenzene	0.289	0.287	0.7	81	-0.01
31		Naphthalene	0.907	0.874	3.6	83	-0.01
32		Benzoic acid	0.198	0.146	26.3#	60	0.01
33		4-Chloroaniline	0.416	0.431	-3.6	91	-0.01
34	C	Hexachlorobutadiene	0.166	0.157	5.4	80	-0.01
35		Caprolactam	0.094	0.093	1.1	74	0.00
36	C	4-Chloro-3-methylphenol	0.313	0.319	-1.9	91	0.00
37		2-Methylnaphthalene	0.651	0.628	3.5	86	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	83	0.00
39		1,2,4,5-Tetrachlorobenzene	0.575	0.546	5.0	79	-0.01
40	P	Hexachlorocyclopentadiene	0.347	0.196	43.5#	46#	-0.01
41	S	2,4,6-Tribromophenol	0.202	0.177	12.4	68	-0.01
42	C	2,4,6-Trichlorophenol	0.430	0.378	12.1	75	-0.01
43		2,4,5-Trichlorophenol	0.412	0.394	4.4	77	-0.01
44	S	2-Fluorobiphenyl	1.317	1.182	10.3	84	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.590	1.508	5.2	85	-0.01
46		2-Chloronaphthalene	1.249	1.163	6.9	86	-0.01
47		2-Nitroaniline	0.412	0.397	3.6	83	-0.01
48		Acenaphthylene	1.845	1.848	-0.2	79	0.00
49		Dimethylphthalate	1.430	1.278	10.6	71	-0.01
50		2,6-Dinitrotoluene	0.314	0.310	1.3	74	0.00
51	C	Acenaphthene	1.237	1.229	0.6	77	0.00
52		3-Nitroaniline	0.389	0.340	12.6	67	-0.01
53	P	2,4-Dinitrophenol	0.093	0.109	-17.2	68	-0.01
54		Dibenzofuran	1.812	1.600	11.7	76	0.00
55	P	4-Nitrophenol	0.282	0.227	19.5	56	-0.01
56		2,4-Dinitrotoluene	0.397	0.431	-8.6	77	0.00
57		Fluorene	1.400	1.263	9.8	80	0.00
58		2,3,4,6-Tetrachlorophenol	0.352	0.309	12.2	69	-0.01
59		Diethylphthalate	1.410	1.267	10.1	73	-0.01
60		4-Chlorophenyl-phenylether	0.564	0.513	9.0	76	-0.01
61		4-Nitroaniline	0.346	0.336	2.9	83	0.00
62		Azobenzene	1.546	1.370	11.4	72	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	80	-0.01
64		4,6-Dinitro-2-methylphenol	0.109	0.090	17.4	63	-0.01
65	c	n-Nitrosodiphenylamine	0.639	0.665	-4.1	82	0.00
66		4-Bromophenyl-phenylether	0.215	0.212	1.4	73	-0.01
67		Hexachlorobenzene	0.242	0.228	5.8	79	0.00
68		Atrazine	0.209	0.180	13.9	61	-0.01
69	C	Pentachlorophenol	0.126	0.106	15.9	59	0.00
70		Phenanthrene	1.046	1.067	-2.0	74	0.00
71		Anthracene	1.026	0.960	6.4	77	-0.01
72		Carbazole	1.003	0.858	14.5	65	-0.01
73		Di-n-butylphthalate	1.111	1.070	3.7	75	-0.01
74	C	Fluoranthene	1.093	0.920	15.8	69	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	60	-0.01
76		Benzidine	0.717	0.676	5.7	56	-0.01
77		Pyrene	1.569	1.554	1.0	68	-0.01
78	S	Terphenyl-d14	0.896	0.882	1.6	69	-0.01
79		Butylbenzylphthalate	0.679	0.697	-2.7	61	-0.01
80		Benzo(a)anthracene	1.174	1.129	3.8	61	-0.01
81		3,3'-Dichlorobenzidine	0.410	0.441	-7.6	64	-0.01
82		Chrysene	1.128	1.147	-1.7	62	0.00
83		Bis(2-ethylhexyl)phthalate	0.858	0.835	2.7	61	-0.01
84	c	Di-n-octyl phthalate	1.449	1.467	-1.2	58	0.00
85		Indeno(1,2,3-cd)pyrene	0.895	1.207	-34.9#	83	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	79	-0.01
87		Benzo(b)fluoranthene	1.308	1.116	14.7	65	-0.01

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88	Benzo(k)fluoranthene	1.070	1.093	-2.1	81	0.00
89 C	Benzo(a)pyrene	1.110	1.101	0.8	75	-0.01
90	Dibenzo(a,h)anthracene	0.955	0.982	-2.8	85	0.00
91	Benzo(g,h,i)perylene	0.989	0.951	3.8	82	-0.01

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