Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012216\
Data File : BF084600.D
Acq On : 22 Jan 2016 20:13
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 23 00:48:07 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	55	-0.08
2		1,4-Dioxane	0.586	0.681	-16.2	61	-0.15
3		Pyridine	1.633	1.398	14.4	43#	-0.19
4		n-Nitrosodimethylamine	0.838	0.830	1.0	51	-0.17
5	S	2-Fluorophenol	1.236	1.183	4.3	55	-0.10
6		Aniline	2.272	2.030	10.7	47#	-0.08
7	S	Phenol-d6	1.553	1.545	0.5	54	-0.08
8		2-Chlorophenol	1.403	1.304	7.1	51	-0.09
9		Benzaldehyde	1.011	0.868	14.1	47#	-0.09
10	C	Phenol	1.766	1.783	-1.0	51	-0.08
11		bis(2-Chloroethyl)ether	1.368	1.300	5.0	54	-0.09
12		1,3-Dichlorobenzene	1.447	1.415	2.2	55	-0.09
13	C	1,4-Dichlorobenzene	1.451	1.371	5.5	49#	-0.09
14		1,2-Dichlorobenzene	1.390	1.476	-6.2	62	-0.08
15		Benzyl Alcohol	1.306	1.158	11.3	46#	-0.08
16		2,2'-oxybis(1-Chloropropane	2.491	2.084	16.3	47#	-0.08
17		2-Methylphenol	1.176	1.171	0.4	51	-0.08
18		Hexachloroethane	0.580	0.578	0.3	52	-0.08
19	P	n-Nitroso-di-n-propylamine	1.051	1.012	3.7	54	-0.08
20		3+4-Methylphenols	1.382	1.434	-3.8	60	-0.08
21	I	Naphthalene-d8	1.000	1.000	0.0	51	-0.08
22		Acetophenone	0.481	0.548	-13.9	54	-0.08
23	S	Nitrobenzene-d5	0.359	0.353	1.7	51	-0.08
24		Nitrobenzene	0.364	0.428	-17.6	54	-0.08
25		Isophorone	0.687	0.699	-1.7	52	-0.08
26	C	2-Nitrophenol	0.166	0.170	-2.4	53	-0.09
27		2,4-Dimethylphenol	0.336	0.316	6.0	45#	-0.07
28		bis(2-Chloroethoxy)methane	0.420	0.438	-4.3	57	-0.08
29	C	2,4-Dichlorophenol	0.280	0.301	-7.5	56	-0.08
30		1,2,4-Trichlorobenzene	0.289	0.317	-9.7	54	-0.08
31		Naphthalene	0.907	0.996	-9.8	57	-0.08
32		Benzoic acid	0.198	0.199	-0.5	49#	-0.05
33		4-Chloroaniline	0.416	0.389	6.5	49#	-0.09
34	C	Hexachlorobutadiene	0.166	0.185	-11.4	57	-0.08
35		Caprolactam	0.094	0.084	10.6	40#	-0.07
36	C	4-Chloro-3-methylphenol	0.313	0.317	-1.3	54	-0.08
37		2-Methylnaphthalene	0.651	0.684	-5.1	56	-0.08
38	I	Acenaphthene-d10	1.000	1.000	0.0	51	-0.08
39		1,2,4,5-Tetrachlorobenzene	0.575	0.606	-5.4	53	-0.08
40	P	Hexachlorocyclopentadiene	0.347	0.205	40.9#	29#	-0.09
41	S	2,4,6-Tribromophenol	0.202	0.191	5.4	45#	-0.09
42	C	2,4,6-Trichlorophenol	0.430	0.436	-1.4	53	-0.08
43		2,4,5-Trichlorophenol	0.412	0.402	2.4	48#	-0.08
44	S	2-Fluorobiphenyl	1.317	1.177	10.6	51	-0.08

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.590	1.881	-18.3	64	-0.08
46		2-Chloronaphthalene	1.249	1.340	-7.3	60	-0.08
47		2-Nitroaniline	0.412	0.433	-5.1	56	-0.09
48		Acenaphthylene	1.845	1.956	-6.0	51	-0.08
49		Dimethylphthalate	1.430	1.517	-6.1	52	-0.08
50		2,6-Dinitrotoluene	0.314	0.312	0.6	46#	-0.08
51	C	Acenaphthene	1.237	1.284	-3.8	49#	-0.08
52		3-Nitroaniline	0.389	0.407	-4.6	49#	-0.08
53	P	2,4-Dinitrophenol	0.093	0.102	-9.7	39#	-0.08
54		Dibenzofuran	1.812	1.698	6.3	49#	-0.08
55	P	4-Nitrophenol	0.282	0.278	1.4	42#	-0.08
56		2,4-Dinitrotoluene	0.397	0.396	0.3	43#	-0.08
57		Fluorene	1.400	1.383	1.2	53	-0.08
58		2,3,4,6-Tetrachlorophenol	0.352	0.365	-3.7	50#	-0.08
59		Diethylphthalate	1.410	1.550	-9.9	55	-0.08
60		4-Chlorophenyl-phenylether	0.564	0.622	-10.3	56	-0.08
61		4-Nitroaniline	0.346	0.388	-12.1	58	-0.07
62		Azobenzene	1.546	1.467	5.1	47#	-0.09
63	I	Phenanthrene-d10	1.000	1.000	0.0	51	-0.09
64		4,6-Dinitro-2-methylphenol	0.109	0.095	12.8	42#	-0.08
65	c	n-Nitrosodiphenylamine	0.639	0.625	2.2	49#	-0.08
66		4-Bromophenyl-phenylether	0.215	0.206	4.2	45#	-0.09
67		Hexachlorobenzene	0.242	0.259	-7.0	57	-0.08
68		Atrazine	0.209	0.229	-9.6	49#	-0.08
69	C	Pentachlorophenol	0.126	0.122	3.2	43#	-0.08
70		Phenanthrene	1.046	1.168	-11.7	52	-0.08
71		Anthracene	1.026	1.009	1.7	52	-0.09
72		Carbazole	1.003	0.872	13.1	42#	-0.09
73		Di-n-butylphthalate	1.111	1.302	-17.2	58	-0.08
74	C	Fluoranthene	1.093	0.959	12.3	46#	-0.09
75	I	Chrysene-d12	1.000	1.000	0.0	45#	-0.09
76		Benzidine	0.717	0.483	32.6#	30#	-0.09
77		Pyrene	1.569	1.327	15.4	43#	-0.09
78	S	Terphenyl-d14	0.896	0.871	2.8	51	-0.09
79		Butylbenzylphthalate	0.679	0.641	5.6	42#	-0.09
80		Benzo(a)anthracene	1.174	1.130	3.7	46#	-0.09
81		3,3'-Dichlorobenzidine	0.410	0.386	5.9	42#	-0.09
82		Chrysene	1.128	0.938	16.8	38#	-0.09
83		Bis(2-ethylhexyl)phthalate	0.858	0.831	3.1	46#	-0.09
84	c	Di-n-octyl phthalate	1.449	1.397	3.6	41#	-0.08
85		Indeno(1,2,3-cd)pyrene	0.895	1.104	-23.4	57	-0.16
86	I	Perylene-d12	1.000	1.000	0.0	51	-0.11
87		Benzo(b)fluoranthene	1.308	1.250	4.4	47#	-0.10

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.070	0.978	8.6	47#	-0.10
89 C	Benzo(a)pyrene	1.110	1.096	1.3	49#	-0.11
90	Dibenzo(a,h)anthracene	0.955	1.016	-6.4	57	-0.16
91	Benzo(g,h,i)perylene	0.989	1.018	-2.9	57	-0.18

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