Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060116\
Data File : BF087622.D
Acq On : 1 Jun 2016 17:58
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 02 08:07:51 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 02 02:13:57 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	103	-0.10
2		1,4-Dioxane	0.601	0.494	17.8	87	-0.09
3		Pyridine	1.313	1.127	14.2	82	-0.11
4		n-Nitrosodimethylamine	1.012	1.052	-4.0	104	-0.10
5	S	2-Fluorophenol	1.138	1.232	-8.3	106	-0.10
6		Aniline	1.990	2.040	-2.5	102	-0.09
7	S	Phenol-d6	1.538	1.558	-1.3	104	-0.07
8		2-Chlorophenol	1.419	1.425	-0.4	103	-0.08
9		Benzaldehyde	0.849	0.787	7.3	104	-0.09
10	C	Phenol	1.679	1.709	-1.8	114	-0.07
11		bis(2-Chloroethyl)ether	1.359	1.333	1.9	102	-0.09
12		1,3-Dichlorobenzene	1.548	1.518	1.9	102	-0.10
13	C	1,4-Dichlorobenzene	1.589	1.558	2.0	102	-0.10
14		1,2-Dichlorobenzene	1.470	1.443	1.8	104	-0.10
15		Benzyl Alcohol	1.121	1.243	-10.9	107	-0.08
16		2,2'-oxybis(1-Chloropropane	3.058	2.830	7.5	98	-0.09
17		2-Methylphenol	1.138	1.154	-1.4	105	-0.08
18		Hexachloroethane	0.593	0.600	-1.2	105	-0.10
19	P	n-Nitroso-di-n-propylamine	1.147	1.153	-0.5	105	-0.08
20		3+4-Methylphenols	1.375	1.546	-12.4	110	-0.07
21	I	Naphthalene-d8	1.000	1.000	0.0	104	-0.09
22		Acetophenone	0.478	0.495	-3.6	107	-0.09
23	S	Nitrobenzene-d5	0.397	0.415	-4.5	105	-0.08
24		Nitrobenzene	0.419	0.431	-2.9	103	-0.09
25		Isophorone	0.712	0.719	-1.0	104	-0.09
26	C	2-Nitrophenol	0.171	0.186	-8.8	106	-0.09
27		2,4-Dimethylphenol	0.297	0.297	0.0	102	-0.08
28		bis(2-Chloroethoxy)methane	0.401	0.386	3.7	101	-0.09
29	C	2,4-Dichlorophenol	0.268	0.280	-4.5	104	-0.07
30		1,2,4-Trichlorobenzene	0.307	0.293	4.6	99	-0.10
31		Naphthalene	1.002	0.998	0.4	106	-0.09
32		Benzoic acid	0.141	0.165	-17.0	112	-0.01
33		4-Chloroaniline	0.369	0.370	-0.3	102	-0.09
34	C	Hexachlorobutadiene	0.186	0.182	2.2	101	-0.10
35		Caprolactam	0.078	0.086	-10.3	110	-0.05
36	C	4-Chloro-3-methylphenol	0.303	0.324	-6.9	106	-0.07
37		2-Methylnaphthalene	0.626	0.608	2.9	102	-0.10
38	I	Acenaphthene-d10	1.000	1.000	0.0	108	-0.10
39		1,2,4,5-Tetrachlorobenzene	0.640	0.605	5.5	102	-0.10
40	P	Hexachlorocyclopentadiene	0.206	0.125	39.3#	58	-0.10
41	S	2,4,6-Tribromophenol	0.192	0.196	-2.1	106	-0.09
42	C	2,4,6-Trichlorophenol	0.370	0.367	0.8	102	-0.08
43		2,4,5-Trichlorophenol	0.376	0.351	6.6	99	-0.07
44	S	2-Fluorobiphenyl	1.419	1.310	7.7	102	-0.10

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.684	1.623	3.6	107	-0.10
46		2-Chloronaphthalene	1.288	1.226	4.8	104	-0.09
47		2-Nitroaniline	0.464	0.493	-6.2	110	-0.08
48		Acenaphthylene	2.040	1.865	8.6	100	-0.10
49		Dimethylphthalate	1.602	1.518	5.2	102	-0.09
50		2,6-Dinitrotoluene	0.323	0.316	2.2	105	-0.08
51	C	Acenaphthene	1.254	1.191	5.0	107	-0.09
52		3-Nitroaniline	0.331	0.339	-2.4	108	-0.07
53	P	2,4-Dinitrophenol	0.138	0.169	-22.5	113	-0.06
54		Dibenzofuran	1.697	1.632	3.8	106	-0.09
55	P	4-Nitrophenol	0.158	0.170	-7.6	113	-0.02
56		2,4-Dinitrotoluene	0.431	0.464	-7.7	110	-0.07
57		Fluorene	1.472	1.364	7.3	101	-0.10
58		2,3,4,6-Tetrachlorophenol	0.289	0.273	5.5	96	-0.09
59		Diethylphthalate	1.558	1.492	4.2	105	-0.09
60		4-Chlorophenyl-phenylether	0.718	0.655	8.8	100	-0.10
61		4-Nitroaniline	0.309	0.346	-12.0	108	-0.07
62		Azobenzene	1.616	1.689	-4.5	108	-0.10
63	I	Phenanthrene-d10	1.000	1.000	0.0	111	-0.09
64		4,6-Dinitro-2-methylphenol	0.127	0.127	0.0	100	-0.06
65	c	n-Nitrosodiphenylamine	0.647	0.615	4.9	103	-0.09
66		4-Bromophenyl-phenylether	0.219	0.209	4.6	105	-0.10
67		Hexachlorobenzene	0.229	0.217	5.2	104	-0.10
68		Atrazine	0.206	0.102	50.5#	54	-0.09
69	C	Pentachlorophenol	0.061	0.053	13.1	85	-0.08
70		Phenanthrene	1.108	1.055	4.8	107	-0.09
71		Anthracene	1.119	1.050	6.2	105	-0.09
72		Carbazole	0.955	0.925	3.1	107	-0.08
73		Di-n-butylphthalate	1.234	1.155	6.4	99	-0.09
74	C	Fluoranthene	1.111	1.057	4.9	104	-0.09
75	I	Chrysene-d12	1.000	1.000	0.0	109	-0.08
76		Benzidine	0.515	0.406	21.2	78	-0.08
77		Pyrene	1.440	1.434	0.4	108	-0.08
78	S	Terphenyl-d14	0.941	0.962	-2.2	109	-0.08
79		Butylbenzylphthalate	0.638	0.680	-6.6	108	-0.08
80		Benzo(a)anthracene	1.138	1.100	3.3	104	-0.07
81		3,3'-Dichlorobenzidine	0.628	0.660	-5.1	117	-0.08
82		Chrysene	1.129	1.047	7.3	103	-0.08
83		Bis(2-ethylhexyl)phthalate	0.932	0.843	9.5	96	-0.09
84	c	Di-n-octyl phthalate	1.421	1.265	11.0	90	-0.09
85		Indeno(1,2,3-cd)pyrene	0.905	0.897	0.9	106	-0.08
86	I	Perylene-d12	1.000	1.000	0.0	96	-0.06
87		Benzo(b)fluoranthene	1.274	1.128	11.5	79	-0.06

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.097	1.263	-15.1	131	-0.07
89 C	Benzo(a)pyrene	1.102	1.117	-1.4	100	-0.06
90	Dibenzo(a,h)anthracene	0.940	1.030	-9.6	104	-0.08
91	Benzo(g,h,i)perylene	0.927	1.023	-10.4	105	-0.09

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