Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF071117\
Data File : BF096661.D
Acq On : 12 Jul 2017 00:03
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 12 01:15:49 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 07 13:02:33 2017
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	66	-0.01
2		1,4-Dioxane	0.643	0.567	11.8	56	-0.05
3		Pyridine	1.763	1.485	15.8	56	-0.06
4		n-Nitrosodimethylamine	0.870	0.772	11.3	58	-0.06
5	S	2-Fluorophenol	1.355	1.244	8.2	62	-0.02
6		Aniline	2.153	2.004	6.9	62	-0.02
7	S	Phenol-d6	1.666	1.603	3.8	64	-0.02
8		2-Chlorophenol	1.445	1.425	1.4	66	-0.01
9		Benzaldehyde	1.043	0.887	15.0	56	-0.02
10	C	Phenol	1.898	1.778	6.3	62	-0.02
11		bis(2-Chloroethyl)ether	1.467	1.352	7.8	61	-0.01
12		1,3-Dichlorobenzene	1.515	1.510	0.3	67	-0.01
13	C	1,4-Dichlorobenzene	1.514	1.523	-0.6	66	-0.01
14		1,2-Dichlorobenzene	1.390	1.387	0.2	68	-0.02
15		Benzyl Alcohol	1.114	1.089	2.2	65	-0.01
16		2,2'-oxybis(1-Chloropropane	2.982	2.853	4.3	63	-0.01
17		2-Methylphenol	1.152	1.115	3.2	64	-0.01
18		Hexachloroethane	0.549	0.506	7.8	60	-0.02
19	P	n-Nitroso-di-n-propylamine	0.992	0.909	8.4	61	-0.02
20		3+4-Methylphenols	1.284	1.323	-3.0	68	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	66	-0.02
22		Acetophenone	0.437	0.428	2.1	67	-0.02
23	S	Nitrobenzene-d5	0.333	0.326	2.1	64	-0.02
24		Nitrobenzene	0.349	0.334	4.3	63	-0.02
25		Isophorone	0.645	0.606	6.0	62	-0.02
26	C	2-Nitrophenol	0.185	0.180	2.7	63	-0.01
27		2,4-Dimethylphenol	0.315	0.302	4.1	64	-0.01
28		bis(2-Chloroethoxy)methane	0.426	0.406	4.7	63	-0.02
29	C	2,4-Dichlorophenol	0.271	0.274	-1.1	67	-0.01
30		1,2,4-Trichlorobenzene	0.256	0.247	3.5	64	-0.01
31		Naphthalene	0.944	0.941	0.3	67	-0.01
32		Benzoic acid	0.264	0.181	31.4#	44#	-0.02
33		4-Chloroaniline	0.392	0.396	-1.0	67	-0.01
34	C	Hexachlorobutadiene	0.131	0.135	-3.1	69	-0.01
35		Caprolactam	0.094	0.092	2.1	67	-0.01
36	C	4-Chloro-3-methylphenol	0.300	0.302	-0.7	68	-0.02
37		2-Methylnaphthalene	0.601	0.593	1.3	67	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	70	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.519	0.521	-0.4	70	-0.02
40	P	Hexachlorocyclopentadiene	0.253	0.129	49.0#	34#	-0.02
41	S	2,4,6-Tribromophenol	0.156	0.157	-0.6	71	-0.02
42	C	2,4,6-Trichlorophenol	0.411	0.389	5.4	65	-0.01
43		2,4,5-Trichlorophenol	0.406	0.397	2.2	68	-0.01
44	S	2-Fluorobiphenyl	1.170	1.190	-1.7	70	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.714	1.696	1.1	70	-0.02
46		2-Chloronaphthalene	1.277	1.259	1.4	69	-0.02
47		2-Nitroaniline	0.465	0.465	0.0	68	-0.01
48		Acenaphthylene	2.024	2.004	1.0	70	-0.02
49		Dimethylphthalate	1.493	1.477	1.1	70	-0.02
50		2,6-Dinitrotoluene	0.330	0.325	1.5	67	-0.01
51	C	Acenaphthene	1.247	1.145	8.2	64	-0.01
52		3-Nitroaniline	0.412	0.421	-2.2	71	-0.01
53	P	2,4-Dinitrophenol	0.175	0.103	41.1#	39#	-0.02
54		Dibenzofuran	1.647	1.640	0.4	69	-0.01
55	P	4-Nitrophenol	0.341	0.298	12.6	59	-0.01
56		2,4-Dinitrotoluene	0.418	0.427	-2.2	68	-0.02
57		Fluorene	1.163	1.215	-4.5	72	-0.01
58		2,3,4,6-Tetrachlorophenol	0.281	0.279	0.7	69	-0.02
59		Diethylphthalate	1.411	1.421	-0.7	71	-0.02
60		4-Chlorophenyl-phenylether	0.506	0.512	-1.2	71	-0.01
61		4-Nitroaniline	0.374	0.401	-7.2	74	-0.01
62		Azobenzene	1.365	1.173	14.1	59	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	69	-0.01
64		4,6-Dinitro-2-methylphenol	0.138	0.101	26.8#	48#	-0.02
65	c	n-Nitrosodiphenylamine	0.702	0.681	3.0	68	-0.02
66		4-Bromophenyl-phenylether	0.195	0.188	3.6	67	-0.02
67		Hexachlorobenzene	0.213	0.213	0.0	70	-0.02
68		Atrazine	0.200	0.203	-1.5	71	-0.01
69	C	Pentachlorophenol	0.126	0.119	5.6	62	-0.01
70		Phenanthrene	1.081	1.063	1.7	70	-0.02
71		Anthracene	1.102	1.094	0.7	70	-0.02
72		Carbazole	1.108	1.072	3.2	68	-0.02
73		Di-n-butylphthalate	1.342	1.341	0.1	70	-0.01
74	C	Fluoranthene	1.060	0.981	7.5	65	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	61	-0.02
76		Benzidine	0.960	0.868	9.6	55	-0.01
77		Pyrene	1.605	1.743	-8.6	67	-0.02
78	S	Terphenyl-d14	0.936	1.065	-13.8	73	-0.02
79		Butylbenzylphthalate	0.813	0.902	-10.9	67	-0.02
80		Benzo(a)anthracene	1.158	1.177	-1.6	64	-0.02
81		3,3'-Dichlorobenzidine	0.476	0.470	1.3	60	-0.02
82		Chrysene	1.213	1.244	-2.6	63	-0.02
83		Bis(2-ethylhexyl)phthalate	0.907	1.078	-18.9	74	-0.02
84	c	Di-n-octyl phthalate	1.786	1.895	-6.1	65	-0.02
85		Indeno(1,2,3-cd)pyrene	0.957	1.132	-18.3	76	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	64	-0.02
87		Benzo(b)fluoranthene	1.253	1.270	-1.4	63	-0.02

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88	Benzo(k)fluoranthene	1.121	1.168	-4.2	69	-0.02
89 C	Benzo(a)pyrene	1.119	1.095	2.1	63	-0.02
90	Dibenzo(a,h)anthracene	0.880	1.036	-17.7	77	-0.04
91	Benzo(g,h,i)perylene	0.912	1.083	-18.7	78	-0.04

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