Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF071017\
Data File : BF096604.D
Acq On : 10 Jul 2017 16:40
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 11 03:10:07 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	65	-0.01
2		1,4-Dioxane	0.643	0.625	2.8	60	-0.04
3		Pyridine	1.763	1.529	13.3	56	-0.04
4		n-Nitrosodimethylamine	0.870	0.813	6.6	59	-0.05
5	S	2-Fluorophenol	1.355	1.316	2.9	63	-0.02
6		Aniline	2.153	2.017	6.3	61	-0.02
7	S	Phenol-d6	1.666	1.599	4.0	62	-0.02
8		2-Chlorophenol	1.445	1.444	0.1	65	-0.01
9		Benzaldehyde	1.043	0.898	13.9	55	-0.02
10	C	Phenol	1.898	1.789	5.7	61	-0.02
11		bis(2-Chloroethyl)ether	1.467	1.365	7.0	60	-0.01
12		1,3-Dichlorobenzene	1.515	1.536	-1.4	67	-0.01
13	C	1,4-Dichlorobenzene	1.514	1.559	-3.0	66	-0.01
14		1,2-Dichlorobenzene	1.390	1.432	-3.0	68	-0.02
15		Benzyl Alcohol	1.114	1.098	1.4	63	-0.01
16		2,2'-oxybis(1-Chloropropane	2.982	2.860	4.1	61	-0.01
17		2-Methylphenol	1.152	1.118	3.0	62	-0.01
18		Hexachloroethane	0.549	0.551	-0.4	64	-0.02
19	P	n-Nitroso-di-n-propylamine	0.992	0.937	5.5	61	-0.02
20		3+4-Methylphenols	1.284	1.327	-3.3	66	-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	0.322	3.3	63	-0.02
24		Nitrobenzene	0.349	0.332	4.9	62	-0.02
25		Isophorone	0.645	0.598	7.3	60	-0.01
26	C	2-Nitrophenol	0.185	0.188	-1.6	65	-0.01
27		2,4-Dimethylphenol	0.315	0.304	3.5	63	-0.01
28		bis(2-Chloroethoxy)methane	0.426	0.399	6.3	61	-0.02
29	C	2,4-Dichlorophenol	0.271	0.279	-3.0	67	-0.02
30		1,2,4-Trichlorobenzene	0.256	0.266	-3.9	68	-0.02
31		Naphthalene	0.944	0.942	0.2	66	-0.01
32		Benzoic acid	0.264	0.251	4.9	60	-0.02
33		4-Chloroaniline	0.392	0.388	1.0	65	-0.01
34	C	Hexachlorobutadiene	0.131	0.138	-5.3	69	-0.01
35		Caprolactam	0.094	0.088	6.4	64	-0.02
36	C	4-Chloro-3-methylphenol	0.300	0.299	0.3	67	-0.02
37		2-Methylnaphthalene	0.601	0.606	-0.8	68	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	68	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.519	0.535	-3.1	71	-0.02
40	P	Hexachlorocyclopentadiene	0.253	0.225	11.1	57	-0.02
41	S	2,4,6-Tribromophenol	0.156	0.178	-14.1	79	-0.02
42	C	2,4,6-Trichlorophenol	0.411	0.407	1.0	67	-0.01
43		2,4,5-Trichlorophenol	0.406	0.411	-1.2	69	-0.02
44	S	2-Fluorobiphenyl	1.170	1.240	-6.0	72	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.714	1.712	0.1	69	-0.02
46		2-Chloronaphthalene	1.277	1.265	0.9	68	-0.02
47		2-Nitroaniline	0.465	0.438	5.8	63	-0.01
48		Acenaphthylene	2.024	2.008	0.8	69	-0.02
49		Dimethylphthalate	1.493	1.482	0.7	69	-0.02
50		2,6-Dinitrotoluene	0.330	0.335	-1.5	68	-0.02
51	C	Acenaphthene	1.247	1.163	6.7	64	-0.02
52		3-Nitroaniline	0.412	0.414	-0.5	69	-0.01
53	P	2,4-Dinitrophenol	0.175	0.172	1.7	64	-0.02
54		Dibenzofuran	1.647	1.646	0.1	69	-0.02
55	P	4-Nitrophenol	0.341	0.319	6.5	62	-0.01
56		2,4-Dinitrotoluene	0.418	0.437	-4.5	69	-0.02
57		Fluorene	1.163	1.246	-7.1	73	-0.01
58		2,3,4,6-Tetrachlorophenol	0.281	0.293	-4.3	71	-0.02
59		Diethylphthalate	1.411	1.435	-1.7	70	-0.02
60		4-Chlorophenyl-phenylether	0.506	0.527	-4.2	72	-0.02
61		4-Nitroaniline	0.374	0.389	-4.0	71	-0.02
62		Azobenzene	1.365	1.278	6.4	64	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	71	-0.01
64		4,6-Dinitro-2-methylphenol	0.138	0.141	-2.2	68	-0.02
65	c	n-Nitrosodiphenylamine	0.702	0.696	0.9	71	-0.02
66		4-Bromophenyl-phenylether	0.195	0.202	-3.6	74	-0.02
67		Hexachlorobenzene	0.213	0.218	-2.3	73	-0.02
68		Atrazine	0.200	0.198	1.0	70	-0.02
69	C	Pentachlorophenol	0.126	0.130	-3.2	68	-0.02
70		Phenanthrene	1.081	1.056	2.3	71	-0.02
71		Anthracene	1.102	1.084	1.6	71	-0.02
72		Carbazole	1.108	1.095	1.2	71	-0.02
73		Di-n-butylphthalate	1.342	1.318	1.8	70	-0.01
74	C	Fluoranthene	1.060	1.063	-0.3	72	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	71	-0.02
76		Benzidine	0.960	0.839	12.6	61	-0.02
77		Pyrene	1.605	1.620	-0.9	72	-0.02
78	S	Terphenyl-d14	0.936	0.971	-3.7	77	-0.02
79		Butylbenzylphthalate	0.813	0.790	2.8	68	-0.02
80		Benzo(a)anthracene	1.158	1.197	-3.4	75	-0.02
81		3,3'-Dichlorobenzidine	0.476	0.481	-1.1	71	-0.02
82		Chrysene	1.213	1.232	-1.6	73	-0.02
83		Bis(2-ethylhexyl)phthalate	0.907	0.915	-0.9	73	-0.02
84	c	Di-n-octyl phthalate	1.786	1.668	6.6	66	-0.02
85		Indeno(1,2,3-cd)pyrene	0.957	0.993	-3.8	77	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	72	-0.03
87		Benzo(b)fluoranthene	1.253	1.198	4.4	67	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.121	1.186	-5.8	79	-0.02
89 C	Benzo(a)pyrene	1.119	1.093	2.3	70	-0.03
90	Dibenzo(a,h)anthracene	0.880	0.934	-6.1	78	-0.04
91	Benzo(g,h,i)perylene	0.912	0.935	-2.5	75	-0.05

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