Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041117\
Data File : BF094378.D
Acq On : 12 Apr 2017 5:32
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 12 13:27:03 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 06 13:20:29 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	71	-0.03
2		1,4-Dioxane	0.569	0.554	2.6	68	-0.08
3		Pyridine	1.550	1.410	9.0	62	-0.06
4		n-Nitrosodimethylamine	0.638	0.598	6.3	64	-0.06
5	S	2-Fluorophenol	1.184	1.045	11.7	63	-0.01
6		Aniline	2.038	1.867	8.4	63	-0.03
7	S	Phenol-d6	1.465	1.319	10.0	63	0.00
8		2-Chlorophenol	1.302	1.209	7.1	64	-0.02
9		Benzaldehyde	0.989	0.868	12.2	62	-0.03
10	C	Phenol	1.667	1.614	3.2	65	0.00
11		bis(2-Chloroethyl)ether	1.303	1.289	1.1	69	-0.03
12		1,3-Dichlorobenzene	1.460	1.484	-1.6	71	-0.03
13	C	1,4-Dichlorobenzene	1.479	1.509	-2.0	71	-0.02
14		1,2-Dichlorobenzene	1.362	1.392	-2.2	72	-0.02
15		Benzyl Alcohol	1.072	0.946	11.8	60	-0.02
16		2,2'-oxybis(1-Chloropropane	2.007	1.710	14.8	59	-0.03
17		2-Methylphenol	1.115	1.061	4.8	66	-0.01
18		Hexachloroethane	0.520	0.443	14.8	58	-0.03
19	P	n-Nitroso-di-n-propylamine	0.941	0.920	2.2	68	-0.03
20		3+4-Methylphenols	1.350	1.383	-2.4	73	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	67	-0.03
22		Acetophenone	0.446	0.477	-7.0	73	-0.03
23	S	Nitrobenzene-d5	0.350	0.360	-2.9	68	-0.02
24		Nitrobenzene	0.346	0.349	-0.9	66	-0.03
25		Isophorone	0.616	0.546	11.4	59	-0.04
26	C	2-Nitrophenol	0.165	0.062	62.4#	23#	-0.03
27		2,4-Dimethylphenol	0.284	0.280	1.4	65	-0.02
28		bis(2-Chloroethoxy)methane	0.406	0.403	0.7	66	-0.03
29	C	2,4-Dichlorophenol	0.266	0.218	18.0	54	-0.02
30		1,2,4-Trichlorobenzene	0.274	0.277	-1.1	67	-0.03
31		Naphthalene	0.962	0.955	0.7	67	-0.03
32		Benzoic acid	0.189	0.002	98.9#	1#	0.00
33		4-Chloroaniline	0.390	0.386	1.0	65	-0.02
34	C	Hexachlorobutadiene	0.140	0.144	-2.9	68	-0.03
35		Caprolactam	0.085	0.059	30.6#	45#	-0.06
36	C	4-Chloro-3-methylphenol	0.277	0.241	13.0	57	-0.01
37		2-Methylnaphthalene	0.641	0.618	3.6	64	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	60	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.547	0.612	-11.9	68	-0.02
40	P	Hexachlorocyclopentadiene	0.256	0.026#	89.8#	6#	-0.03
41	S	2,4,6-Tribromophenol	0.158	0.086	45.6#	32#	-0.03
42	C	2,4,6-Trichlorophenol	0.387	0.243	37.2#	37#	-0.02
43		2,4,5-Trichlorophenol	0.419	0.277	33.9#	38#	0.00
44	S	2-Fluorobiphenyl	1.311	1.452	-10.8	67	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.613	1.697	-5.2	63	-0.03
46		2-Chloronaphthalene	1.263	1.303	-3.2	63	-0.02
47		2-Nitroaniline	0.375	0.382	-1.9	59	-0.02
48		Acenaphthylene	2.099	2.070	1.4	59	-0.03
49		Dimethylphthalate	1.422	1.518	-6.8	65	-0.03
50		2,6-Dinitrotoluene	0.326	0.325	0.3	58	-0.02
51	C	Acenaphthene	1.225	1.196	2.4	60	-0.03
52		3-Nitroaniline	0.383	0.407	-6.3	63	-0.02
53	P	2,4-Dinitrophenol	0.155	0.000#	100.0#	0#	-9.59#
54		Dibenzofuran	1.667	1.597	4.2	57	-0.03
55	P	4-Nitrophenol	0.300	0.110	63.3#	21#	0.00
56		2,4-Dinitrotoluene	0.409	0.363	11.2	51	-0.03
57		Fluorene	1.382	1.325	4.1	62	-0.03
58		2,3,4,6-Tetrachlorophenol	0.287	0.106	63.1#	22#	-0.02
59		Diethylphthalate	1.405	1.418	-0.9	61	-0.03
60		4-Chlorophenyl-phenylether	0.589	0.584	0.8	63	-0.02
61		4-Nitroaniline	0.367	0.386	-5.2	61	-0.03
62		Azobenzene	1.329	1.201	9.6	54	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	55	-0.03
64		4,6-Dinitro-2-methylphenol	0.122	0.001	99.2#	0#	-0.03
65	c	n-Nitrosodiphenylamine	0.692	0.732	-5.8	59	-0.03
66		4-Bromophenyl-phenylether	0.197	0.211	-7.1	59	-0.03
67		Hexachlorobenzene	0.199	0.201	-1.0	56	-0.02
68		Atrazine	0.207	0.216	-4.3	57	-0.02
69	C	Pentachlorophenol	0.124	0.015	87.9#	6#	-0.01
70		Phenanthrene	1.113	1.117	-0.4	56	-0.02
71		Anthracene	1.146	1.154	-0.7	56	-0.02
72		Carbazole	1.175	1.170	0.4	55	-0.02
73		Di-n-butylphthalate	1.222	1.297	-6.1	58	-0.02
74	C	Fluoranthene	1.139	1.168	-2.5	57	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	60	-0.02
76		Benzidine	0.792	0.730	7.8	54	-0.02
77		Pyrene	1.451	1.383	4.7	57	-0.02
78	S	Terphenyl-d14	0.892	0.886	0.7	60	-0.02
79		Butylbenzylphthalate	0.618	0.620	-0.3	57	-0.02
80		Benzo(a)anthracene	1.166	1.154	1.0	59	-0.02
81		3,3'-Dichlorobenzidine	0.417	0.424	-1.7	58	-0.02
82		Chrysene	1.082	1.078	0.4	60	-0.02
83		Bis(2-ethylhexyl)phthalate	0.844	0.870	-3.1	59	-0.02
84	c	Di-n-octyl phthalate	1.410	1.480	-5.0	59	-0.02
85		Indeno(1,2,3-cd)pyrene	0.937	0.812	13.3	53	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	56	-0.02
87		Benzo(b)fluoranthene	1.226	1.353	-10.4	60	-0.02

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88	Benzo(k)fluoranthene	1.199	1.177	1.8	53	-0.02
89 C	Benzo(a)pyrene	1.129	1.141	-1.1	55	-0.02
90	Dibenzo(a,h)anthracene	0.956	0.926	3.1	54	-0.02
91	Benzo(g,h,i)perylene	0.964	0.893	7.4	52	-0.03

(#) = Out of Range SPCC's out = 2 CCC's out = 3