Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071315\
Data File : BF080445.D
Acq On : 14 Jul 2015 00:25
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 14 08:31:19 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 14 00:52:37 2015
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	-0.02
2		1,4-Dioxane	0.562	0.601	-6.9	68	-0.07
3		Pyridine	1.248	1.267	-1.5	67	-0.10
4		n-Nitrosodimethylamine	0.684	0.714	-4.4	67	-0.06
5	S	2-Fluorophenol	1.167	1.260	-8.0	68	-0.02
6		Aniline	1.939	2.110	-8.8	72	-0.01
7	S	Phenol-d6	1.555	1.510	2.9	63	-0.02
8		2-Chlorophenol	1.406	1.360	3.3	60	-0.01
9		Benzaldehyde	0.860	0.837	2.7	63	-0.01
10	C	Phenol	1.718	1.775	-3.3	68	-0.01
11		bis(2-Chloroethyl)ether	1.415	1.322	6.6	61	-0.01
12		1,3-Dichlorobenzene	1.590	1.707	-7.4	72	-0.01
13	C	1,4-Dichlorobenzene	1.793	1.862	-3.8	68	-0.01
14		1,2-Dichlorobenzene	1.699	1.692	0.4	66	-0.01
15		Benzyl Alcohol	1.280	1.149	10.2	58	-0.01
16		2,2'-oxybis(1-Chloropropane	2.040	2.379	-16.6	76	-0.01
17		2-Methylphenol	1.222	1.083	11.4	54	-0.01
18		Hexachloroethane	0.553	0.562	-1.6	64	-0.01
19	P	n-Nitroso-di-n-propylamine	1.064	1.005	5.5	63	-0.02
20		3+4-Methylphenols	1.536	1.605	-4.5	65	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	66	-0.01
22		Acetophenone	0.479	0.481	-0.4	61	-0.02
23	S	Nitrobenzene-d5	0.351	0.383	-9.1	71	-0.01
24		Nitrobenzene	0.388	0.373	3.9	57	-0.01
25		Isophorone	0.628	0.703	-11.9	72	-0.02
26	C	2-Nitrophenol	0.186	0.168	9.7	55	-0.01
27		2,4-Dimethylphenol	0.293	0.321	-9.6	70	-0.01
28		bis(2-Chloroethoxy)methane	0.380	0.420	-10.5	69	-0.02
29	C	2,4-Dichlorophenol	0.271	0.299	-10.3	68	-0.02
30		1,2,4-Trichlorobenzene	0.358	0.342	4.5	60	-0.01
31		Naphthalene	1.094	1.139	-4.1	69	-0.01
32		Benzoic acid	0.206	0.180	12.6	57	-0.03
33		4-Chloroaniline	0.371	0.411	-10.8	67	-0.01
34	C	Hexachlorobutadiene	0.208	0.209	-0.5	65	-0.01
35		Caprolactam	0.077	0.068	11.7	58	-0.04
36	C	4-Chloro-3-methylphenol	0.287	0.316	-10.1	68	-0.02
37		2-Methylnaphthalene	0.722	0.671	7.1	59	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	63	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.669	0.681	-1.8	62	-0.01
40	P	Hexachlorocyclopentadiene	0.349	0.334	4.3	58	-0.01
41	S	2,4,6-Tribromophenol	0.195	0.159	18.5	46#	-0.02
42	C	2,4,6-Trichlorophenol	0.385	0.400	-3.9	60	-0.02
43		2,4,5-Trichlorophenol	0.414	0.458	-10.6	65	-0.01
44	S	2-Fluorobiphenyl	1.485	1.546	-4.1	65	-0.01

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 14 00:52:37 2015
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)
45		1,1'-Biphenyl	1.743	1.742	0.1	61	-0.01
46		2-Chloronaphthalene	1.287	1.246	3.2	58	-0.02
47		2-Nitroaniline	0.380	0.320	15.8	50#	-0.01
48		Acenaphthylene	2.201	2.350	-6.8	65	-0.01
49		Dimethylphthalate	1.512	1.525	-0.9	62	-0.02
50		2,6-Dinitrotoluene	0.339	0.289	14.7	51	-0.01
51	C	Acenaphthene	1.399	1.423	-1.7	62	-0.01
52		3-Nitroaniline	0.331	0.292	11.8	52	-0.01
53	P	2,4-Dinitrophenol	0.128	0.134	-4.7	60	-0.02
54		Dibenzofuran	1.853	1.951	-5.3	65	-0.01
55	P	4-Nitrophenol	0.237	0.221	6.8	56	-0.01
56		2,4-Dinitrotoluene	0.397	0.426	-7.3	65	-0.01
57		Fluorene	1.561	1.583	-1.4	64	-0.01
58		2,3,4,6-Tetrachlorophenol	0.363	0.344	5.2	54	-0.01
59		Diethylphthalate	1.480	1.379	6.8	59	-0.01
60		4-Chlorophenyl-phenylether	0.752	0.691	8.1	56	-0.01
61		4-Nitroaniline	0.320	0.293	8.4	55	-0.02
62		Azobenzene	1.598	1.478	7.5	56	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	51	-0.01
64		4,6-Dinitro-2-methylphenol	0.105	0.121	-15.2	55	-0.01
65	c	n-Nitrosodiphenylamine	0.623	0.751	-20.5#	59	-0.02
66		4-Bromophenyl-phenylether	0.213	0.234	-9.9	55	-0.01
67		Hexachlorobenzene	0.214	0.234	-9.3	55	-0.02
68		Atrazine	0.194	0.213	-9.8	49#	-0.02
69	C	Pentachlorophenol	0.122	0.124	-1.6	48#	-0.01
70		Phenanthrene	1.202	1.258	-4.7	54	-0.01
71		Anthracene	1.113	1.325	-19.0	59	-0.02
72		Carbazole	1.027	1.223	-19.1	59	-0.02
73		Di-n-butylphthalate	1.086	1.327	-22.2	55	-0.02
74	C	Fluoranthene	1.246	1.307	-4.9	53	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	55	-0.06
76		Benzidine	0.391	0.511	-30.7#	70	-0.03
77		Pyrene	1.281	1.333	-4.1	58	-0.02
78	S	Terphenyl-d14	0.921	0.938	-1.8	56	-0.03
79		Butylbenzylphthalate	0.486	0.453	6.8	48#	-0.03
80		Benzo(a)anthracene	1.260	1.260	0.0	54	-0.06
81		3,3'-Dichlorobenzidine	0.377	0.400	-6.1	55	-0.05
82		Chrysene	1.161	1.174	-1.1	54	-0.06
83		Bis(2-ethylhexyl)phthalate	0.762	0.688	9.7	48#	-0.06
84	c	Di-n-octyl phthalate	1.343	1.147	14.6	45#	-0.07
85		Indeno(1,2,3-cd)pyrene	1.442	2.443	-69.4#	94	-0.11
86	I	Perylene-d12	1.000	1.000	0.0	66	-0.08
87		Benzo(b)fluoranthene	1.389	1.247	10.2	60	-0.08

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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)
88	Benzo(k)fluoranthene	1.136	1.019	10.3	60	-0.08
89 C	Benzo(a)pyrene	1.217	1.158	4.8	64	-0.09
90	Dibenzo(a,h)anthracene	1.206	1.603	-32.9#	91	-0.11
91	Benzo(g,h,i)perylene	1.198	1.783	-48.8#	101	-0.10

(#) = Out of Range SPCC's out = 0 CCC's out = 1