Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071315\
Data File : BF080447.D
Acq On : 14 Jul 2015 4:14
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 14 07:20:39 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	77	-0.02
2		1,4-Dioxane	0.562	0.539	4.1	71	-0.07
3		Pyridine	1.248	1.300	-4.2	80	-0.13
4		n-Nitrosodimethylamine	0.684	0.777	-13.6	84	-0.07
5	S	2-Fluorophenol	1.167	1.242	-6.4	78	-0.02
6		Aniline	1.939	2.254	-16.2	90	-0.01
7	S	Phenol-d6	1.555	1.627	-4.6	80	-0.02
8		2-Chlorophenol	1.406	1.368	2.7	71	-0.01
9		Benzaldehyde	0.860	0.857	0.3	75	-0.01
10	C	Phenol	1.718	1.845	-7.4	82	-0.01
11		bis(2-Chloroethyl)ether	1.415	1.347	4.8	72	-0.01
12		1,3-Dichlorobenzene	1.590	1.711	-7.6	84	-0.01
13	C	1,4-Dichlorobenzene	1.793	1.925	-7.4	82	-0.01
14		1,2-Dichlorobenzene	1.699	1.676	1.4	76	-0.01
15		Benzyl Alcohol	1.280	1.253	2.1	74	-0.01
16		2,2'-oxybis(1-Chloropropane	2.040	2.393	-17.3	88	-0.01
17		2-Methylphenol	1.222	1.146	6.2	67	-0.01
18		Hexachloroethane	0.553	0.584	-5.6	78	-0.01
19	P	n-Nitroso-di-n-propylamine	1.064	1.150	-8.1	84	-0.02
20		3+4-Methylphenols	1.536	1.652	-7.6	78	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	80	-0.01
22		Acetophenone	0.479	0.473	1.3	73	-0.02
23	S	Nitrobenzene-d5	0.351	0.394	-12.3	88	-0.01
24		Nitrobenzene	0.388	0.392	-1.0	73	-0.01
25		Isophorone	0.628	0.719	-14.5	89	-0.02
26	C	2-Nitrophenol	0.186	0.175	5.9	70	-0.01
27		2,4-Dimethylphenol	0.293	0.328	-11.9	87	-0.01
28		bis(2-Chloroethoxy)methane	0.380	0.416	-9.5	83	-0.02
29	C	2,4-Dichlorophenol	0.271	0.305	-12.5	84	-0.02
30		1,2,4-Trichlorobenzene	0.358	0.355	0.8	76	-0.01
31		Naphthalene	1.094	1.130	-3.3	83	-0.01
32		Benzoic acid	0.206	0.200	2.9	77	-0.03
33		4-Chloroaniline	0.371	0.436	-17.5	87	-0.01
34	C	Hexachlorobutadiene	0.208	0.204	1.9	77	-0.01
35		Caprolactam	0.077	0.082	-6.5	85	-0.04
36	C	4-Chloro-3-methylphenol	0.287	0.330	-15.0	86	-0.02
37		2-Methylnaphthalene	0.722	0.694	3.9	74	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	79	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.669	0.660	1.3	75	-0.01
40	P	Hexachlorocyclopentadiene	0.349	0.332	4.9	73	-0.01
41	S	2,4,6-Tribromophenol	0.195	0.173	11.3	63	-0.02
42	C	2,4,6-Trichlorophenol	0.385	0.395	-2.6	75	-0.02
43		2,4,5-Trichlorophenol	0.414	0.450	-8.7	81	-0.01
44	S	2-Fluorobiphenyl	1.485	1.511	-1.8	80	-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	1.760	-1.0	78	-0.01
46		2-Chloronaphthalene	1.287	1.253	2.6	74	-0.02
47		2-Nitroaniline	0.380	0.373	1.8	73	-0.01
48		Acenaphthylene	2.201	2.333	-6.0	82	-0.01
49		Dimethylphthalate	1.512	1.669	-10.4	85	-0.02
50		2,6-Dinitrotoluene	0.339	0.323	4.7	72	-0.01
51	C	Acenaphthene	1.399	1.478	-5.6	82	-0.01
52		3-Nitroaniline	0.331	0.370	-11.8	83	-0.01
53	P	2,4-Dinitrophenol	0.128	0.169	-32.0#	96	-0.02
54		Dibenzofuran	1.853	1.994	-7.6	85	-0.01
55	P	4-Nitrophenol	0.237	0.265	-11.8	85	-0.01
56		2,4-Dinitrotoluene	0.397	0.483	-21.7	93	-0.01
57		Fluorene	1.561	1.644	-5.3	83	-0.01
58		2,3,4,6-Tetrachlorophenol	0.363	0.390	-7.4	77	-0.01
59		Diethylphthalate	1.480	1.602	-8.2	87	-0.01
60		4-Chlorophenyl-phenylether	0.752	0.752	0.0	77	-0.01
61		4-Nitroaniline	0.320	0.358	-11.9	85	-0.01
62		Azobenzene	1.598	1.656	-3.6	79	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	74	-0.01
64		4,6-Dinitro-2-methylphenol	0.105	0.129	-22.9	87	-0.01
65	c	n-Nitrosodiphenylamine	0.623	0.671	-7.7	77	-0.02
66		4-Bromophenyl-phenylether	0.213	0.224	-5.2	76	-0.01
67		Hexachlorobenzene	0.214	0.215	-0.5	75	-0.02
68		Atrazine	0.194	0.236	-21.6	79	-0.02
69	C	Pentachlorophenol	0.122	0.132	-8.2	74	-0.01
70		Phenanthrene	1.202	1.267	-5.4	80	-0.01
71		Anthracene	1.113	1.252	-12.5	81	-0.02
72		Carbazole	1.027	1.179	-14.8	83	-0.02
73		Di-n-butylphthalate	1.086	1.261	-16.1	77	-0.02
74	C	Fluoranthene	1.246	1.405	-12.8	82	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	100	-0.05
76		Benzidine	0.391	0.503	-28.6#	125	-0.02
77		Pyrene	1.281	1.096	14.4	86	-0.02
78	S	Terphenyl-d14	0.921	0.760	17.5	83	-0.03
79		Butylbenzylphthalate	0.486	0.454	6.6	88	-0.03
80		Benzo(a)anthracene	1.260	1.241	1.5	97	-0.05
81		3,3'-Dichlorobenzidine	0.377	0.404	-7.2	101	-0.05
82		Chrysene	1.161	1.170	-0.8	98	-0.05
83		Bis(2-ethylhexyl)phthalate	0.762	0.680	10.8	85	-0.06
84	c	Di-n-octyl phthalate	1.343	1.256	6.5	89	-0.06
85		Indeno(1,2,3-cd)pyrene	1.442	2.153	-49.3#	151#	-0.09
86	I	Perylene-d12	1.000	1.000	0.0	120	-0.08
87		Benzo(b)fluoranthene	1.389	1.175	15.4	102	-0.07

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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	1.244	-9.5	133	-0.07
89 C	Benzo(a)pyrene	1.217	1.213	0.3	122	-0.08
90	Dibenzo(a,h)anthracene	1.206	1.458	-20.9	150#	-0.10
91	Benzo(g,h,i)perylene	1.198	1.538	-28.4#	159#	-0.09

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