Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061015\
Data File : BF079879.D
Acq On : 11 Jun 2015 00:06
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 11 11:42:53 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 15:11:40 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	114	0.00
2		1,4-Dioxane	0.555	0.505	9.0	103	-0.02
3		Pyridine	1.490	1.496	-0.4	107	-0.03
4		n-Nitrosodimethylamine	0.746	0.621	16.8	91	-0.01
5	S	2-Fluorophenol	1.214	1.147	5.5	105	0.00
6		Aniline	2.262	2.165	4.3	105	0.00
7	S	Phenol-d6	1.570	1.557	0.8	107	0.00
8		2-Chlorophenol	1.368	1.317	3.7	112	0.00
9		Benzaldehyde	0.921	0.873	5.2	103	-0.01
10	C	Phenol	1.673	1.685	-0.7	110	0.00
11		bis(2-Chloroethyl)ether	1.306	1.283	1.8	115	0.00
12		1,3-Dichlorobenzene	1.568	1.541	1.7	111	0.00
13	C	1,4-Dichlorobenzene	1.777	1.589	10.6	101	0.00
14		1,2-Dichlorobenzene	1.543	1.434	7.1	102	-0.01
15		Benzyl Alcohol	1.266	1.207	4.7	100	0.00
16		2,2'-oxybis(1-Chloropropane	2.025	1.950	3.7	110	0.00
17		2-Methylphenol	1.177	1.160	1.4	106	0.00
18		Hexachloroethane	0.552	0.530	4.0	111	0.00
19	P	n-Nitroso-di-n-propylamine	1.094	1.020	6.8	112	0.00
20		3+4-Methylphenols	1.521	1.481	2.6	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	114	0.00
22		Acetophenone	0.496	0.485	2.2	109	0.00
23	S	Nitrobenzene-d5	0.346	0.330	4.6	99	0.00
24		Nitrobenzene	0.348	0.328	5.7	107	0.00
25		Isophorone	0.722	0.705	2.4	113	0.00
26	C	2-Nitrophenol	0.139	0.111	20.1#	87	0.00
27		2,4-Dimethylphenol	0.326	0.316	3.1	112	0.00
28		bis(2-Chloroethoxy)methane	0.437	0.393	10.1	99	0.00
29	C	2,4-Dichlorophenol	0.278	0.262	5.8	104	0.00
30		1,2,4-Trichlorobenzene	0.357	0.329	7.8	102	0.00
31		Naphthalene	1.065	1.055	0.9	115	0.00
32		Benzoic acid	0.098	0.077	21.4	83	0.02
33		4-Chloroaniline	0.430	0.438	-1.9	115	0.00
34	C	Hexachlorobutadiene	0.196	0.187	4.6	103	-0.01
35		Caprolactam	0.081	0.084	-3.7	113	0.02
36	C	4-Chloro-3-methylphenol	0.301	0.301	0.0	107	0.00
37		2-Methylnaphthalene	0.757	0.727	4.0	111	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
39		1,2,4,5-Tetrachlorobenzene	0.699	0.663	5.2	98	0.00
40	P	Hexachlorocyclopentadiene	0.334	0.268	19.8	87	0.00
41	S	2,4,6-Tribromophenol	0.173	0.178	-2.9	112	0.00
42	C	2,4,6-Trichlorophenol	0.422	0.401	5.0	98	0.00
43		2,4,5-Trichlorophenol	0.436	0.483	-10.8	118	0.00
44	S	2-Fluorobiphenyl	1.422	1.424	-0.1	108	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.658	1.639	1.1	104	0.00
46		2-Chloronaphthalene	1.354	1.406	-3.8	112	0.00
47		2-Nitroaniline	0.297	0.294	1.0	109	0.00
48		Acenaphthylene	2.302	2.356	-2.3	112	0.00
49		Dimethylphthalate	1.559	1.593	-2.2	114	0.00
50		2,6-Dinitrotoluene	0.270	0.290	-7.4	117	0.00
51	C	Acenaphthene	1.347	1.329	1.3	106	0.00
52		3-Nitroaniline	0.333	0.339	-1.8	111	0.00
53	P	2,4-Dinitrophenol	0.064	0.040#	37.5#	63	0.00
54		Dibenzofuran	1.886	1.899	-0.7	107	0.00
55	P	4-Nitrophenol	0.242	0.232	4.1	103	0.00
56		2,4-Dinitrotoluene	0.347	0.308	11.2	90	0.00
57		Fluorene	1.670	1.673	-0.2	106	0.00
58		2,3,4,6-Tetrachlorophenol	0.322	0.352	-9.3	113	0.00
59		Diethylphthalate	1.527	1.620	-6.1	114	0.00
60		4-Chlorophenyl-phenylether	0.723	0.754	-4.3	115	0.00
61		4-Nitroaniline	0.326	0.360	-10.4	114	0.00
62		Azobenzene	1.503	1.594	-6.1	114	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	111	-0.01
64		4,6-Dinitro-2-methylphenol	0.057	0.042	26.3#	80	0.00
65	c	n-Nitrosodiphenylamine	0.664	0.696	-4.8	123	0.00
66		4-Bromophenyl-phenylether	0.231	0.219	5.2	113	0.00
67		Hexachlorobenzene	0.243	0.254	-4.5	113	0.00
68		Atrazine	0.188	0.183	2.7	106	-0.01
69	C	Pentachlorophenol	0.089	0.081	9.0	101	-0.01
70		Phenanthrene	1.195	1.176	1.6	113	0.00
71		Anthracene	1.170	1.150	1.7	105	-0.01
72		Carbazole	1.083	1.063	1.8	110	-0.01
73		Di-n-butylphthalate	1.124	1.222	-8.7	116	-0.01
74	C	Fluoranthene	1.253	1.296	-3.4	112	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	111	-0.11
76		Benzidine	0.596	0.609	-2.2	104	-0.03
77		Pyrene	1.246	1.277	-2.5	111	-0.03
78	S	Terphenyl-d14	0.887	0.862	2.8	106	-0.05
79		Butylbenzylphthalate	0.417	0.434	-4.1	113	-0.06
80		Benzo(a)anthracene	1.189	1.200	-0.9	109	-0.10
81		3,3'-Dichlorobenzidine	0.371	0.412	-11.1	117	-0.10
82		Chrysene	1.116	1.159	-3.9	111	-0.10
83		Bis(2-ethylhexyl)phthalate	0.645	0.699	-8.4	119	-0.11
84	c	Di-n-octyl phthalate	1.002	1.110	-10.8	121	-0.19
85		Indeno(1,2,3-cd)pyrene	1.178	1.324	-12.4	119	-0.21
86	I	Perylene-d12	1.000	1.000	0.0	117	-0.21
87		Benzo(b)fluoranthene	1.190	1.313	-10.3	128	-0.18

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.245	1.145	8.0	99	-0.19
89 C	Benzo(a)pyrene	1.125	1.155	-2.7	115	-0.18
90	Dibenzo(a,h)anthracene	1.042	1.116	-7.1	116	-0.21
91	Benzo(g,h,i)perylene	1.100	1.128	-2.5	114	-0.21

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