Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079402.D
Acq On : 21 May 2015 14:25
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 22 01:14:42 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 22 01:11:00 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	59	-0.06
2		1,4-Dioxane	0.505	0.744	-47.3#	87	-0.08
3		Pyridine	1.478	1.659	-12.2	70	-0.09
4		n-Nitrosodimethylamine	0.633	0.703	-11.1	67	-0.08
5	S	2-Fluorophenol	1.162	1.253	-7.8	65	-0.07
6		Aniline	2.168	2.289	-5.6	63	-0.06
7	S	Phenol-d6	1.536	1.645	-7.1	68	-0.06
8		2-Chlorophenol	1.318	1.395	-5.8	67	-0.07
9		Benzaldehyde	1.035	0.967	6.6	57	-0.06
10	C	Phenol	1.782	1.899	-6.6	64	-0.06
11		bis(2-Chloroethyl)ether	1.306	1.382	-5.8	68	-0.06
12		1,3-Dichlorobenzene	1.481	1.576	-6.4	68	-0.06
13	C	1,4-Dichlorobenzene	1.524	1.620	-6.3	66	-0.06
14		1,2-Dichlorobenzene	1.419	1.543	-8.7	67	-0.06
15		Benzyl Alcohol	1.140	1.167	-2.4	64	-0.05
16		2,2'-oxybis(1-Chloropropane	1.998	2.000	-0.1	63	-0.06
17		2-Methylphenol	1.060	1.206	-13.8	71	-0.06
18		Hexachloroethane	0.525	0.558	-6.3	64	-0.06
19	P	n-Nitroso-di-n-propylamine	1.037	1.114	-7.4	64	-0.06
20		3+4-Methylphenols	1.413	1.566	-10.8	67	-0.06
21	I	Naphthalene-d8	1.000	1.000	0.0	60	-0.06
22		Acetophenone	0.452	0.478	-5.8	68	-0.06
23	S	Nitrobenzene-d5	0.348	0.378	-8.6	68	-0.06
24		Nitrobenzene	0.345	0.373	-8.1	71	-0.06
25		Isophorone	0.645	0.685	-6.2	66	-0.06
26	C	2-Nitrophenol	0.143	0.150	-4.9	63	-0.06
27		2,4-Dimethylphenol	0.286	0.308	-7.7	66	-0.05
28		bis(2-Chloroethoxy)methane	0.398	0.426	-7.0	65	-0.06
29	C	2,4-Dichlorophenol	0.254	0.279	-9.8	69	-0.06
30		1,2,4-Trichlorobenzene	0.279	0.289	-3.6	65	-0.06
31		Naphthalene	0.953	1.018	-6.8	68	-0.06
32		Benzoic acid	0.164	0.194	-18.3	68	-0.06
33		4-Chloroaniline	0.403	0.431	-6.9	69	-0.06
34	C	Hexachlorobutadiene	0.161	0.164	-1.9	66	-0.06
35		Caprolactam	0.082	0.095	-15.9	71	-0.07
36	C	4-Chloro-3-methylphenol	0.267	0.307	-15.0	68	-0.06
37		2-Methylnaphthalene	0.660	0.688	-4.2	65	-0.06
38	I	Acenaphthene-d10	1.000	1.000	0.0	60	-0.06
39		1,2,4,5-Tetrachlorobenzene	0.563	0.571	-1.4	65	-0.07
40	P	Hexachlorocyclopentadiene	0.283	0.229	19.1	50#	-0.06
41	S	2,4,6-Tribromophenol	0.216	0.232	-7.4	64	-0.06
42	C	2,4,6-Trichlorophenol	0.387	0.411	-6.2	65	-0.06
43		2,4,5-Trichlorophenol	0.392	0.429	-9.4	67	-0.06
44	S	2-Fluorobiphenyl	1.436	1.581	-10.1	69	-0.06
45		1,1'-Biphenyl	1.590	1.669	-5.0	67	-0.06
46		2-Chloronaphthalene	1.240	1.332	-7.4	70	-0.06
47		2-Nitroaniline	0.359	0.389	-8.4	65	-0.06
48		Acenaphthylene	2.036	2.168	-6.5	68	-0.07

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49		Dimethylphthalate	1.468	1.611	-9.7	71	-0.06
50		2,6-Dinitrotoluene	0.290	0.323	-11.4	69	-0.06
51	C	Acenaphthene	1.214	1.236	-1.8	64	-0.06
52		3-Nitroaniline	0.362	0.420	-16.0	72	-0.06
53	P	2,4-Dinitrophenol	0.085	0.067	21.2	49#	-0.06
54		Dibenzofuran	1.754	1.904	-8.6	68	-0.06
55	P	4-Nitrophenol	0.286	0.262	8.4	57	-0.05
56		2,4-Dinitrotoluene	0.383	0.430	-12.3	66	-0.06
57		Fluorene	1.440	1.552	-7.8	69	-0.06
58		2,3,4,6-Tetrachlorophenol	0.320	0.331	-3.4	63	-0.06
59		Diethylphthalate	1.454	1.562	-7.4	68	-0.06
60		4-Chlorophenyl-phenylether	0.639	0.674	-5.5	66	-0.06
61		4-Nitroaniline	0.362	0.422	-16.6	70	-0.06
62		Azobenzene	1.433	1.487	-3.8	64	-0.06
63	I	Phenanthrene-d10	1.000	1.000	0.0	61	-0.07
64		4,6-Dinitro-2-methylphenol	0.080	0.074	7.5	57	-0.06
65	c	n-Nitrosodiphenylamine	0.666	0.706	-6.0	66	-0.06
66		4-Bromophenyl-phenylether	0.208	0.220	-5.8	69	-0.06
67		Hexachlorobenzene	0.238	0.248	-4.2	68	-0.07
68		Atrazine	0.194	0.200	-3.1	67	-0.07
69	C	Pentachlorophenol	0.120	0.105	12.5	54	-0.07
70		Phenanthrene	1.119	1.196	-6.9	68	-0.06
71		Anthracene	1.098	1.208	-10.0	70	-0.06
72		Carbazole	1.046	1.157	-10.6	70	-0.06
73		Di-n-butylphthalate	1.255	1.401	-11.6	68	-0.06
74	C	Fluoranthene	1.166	1.306	-12.0	71	-0.07
75	I	Chrysene-d12	1.000	1.000	0.0	58	-0.06
76		Benzidine	0.539	0.585	-8.5	62	-0.07
77		Pyrene	1.152	1.268	-10.1	68	-0.07
78	S	Terphenyl-d14	0.835	0.897	-7.4	67	-0.07
79		Butylbenzylphthalate	0.504	0.613	-21.6	70	-0.06
80		Benzo(a)anthracene	1.095	1.199	-9.5	67	-0.06
81		3,3'-Dichlorobenzidine	0.390	0.447	-14.6	68	-0.06
82		Chrysene	1.053	1.147	-8.9	69	-0.06
83		Bis(2-ethylhexyl)phthalate	0.763	0.902	-18.2	68	-0.05
84	c	Di-n-octyl phthalate	1.344	1.599	-19.0	73	-0.03
85		Indeno(1,2,3-cd)pyrene	1.342	1.532	-14.2	70	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	64	-0.02
87		Benzo(b)fluoranthene	1.095	1.172	-7.0	70	-0.03
88		Benzo(k)fluoranthene	1.060	1.094	-3.2	68	-0.03
89	C	Benzo(a)pyrene	1.008	1.102	-9.3	72	-0.02
90		Dibenzo(a,h)anthracene	1.071	1.183	-10.5	73	-0.05
91		Benzo(g,h,i)perylene	1.074	1.182	-10.1	73	-0.05

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