Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079420.D
Acq On : 21 May 2015 23:27
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: May 22 03:34:48 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 03:34:28 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	63	-0.06
2		1,4-Dioxane	0.505	0.629	-24.6	79	-0.08
3		Pyridine	1.478	1.347	8.9	61	-0.09
4		n-Nitrosodimethylamine	0.633	0.664	-4.9	68	-0.08
5	S	2-Fluorophenol	1.162	1.248	-7.4	69	-0.07
6		Aniline	2.168	2.260	-4.2	66	-0.06
7	S	Phenol-d6	1.536	1.597	-4.0	70	-0.07
8		2-Chlorophenol	1.318	1.440	-9.3	74	-0.07
9		Benzaldehyde	1.035	0.956	7.6	60	-0.06
10	C	Phenol	1.782	1.900	-6.6	68	-0.06
11		bis(2-Chloroethyl)ether	1.306	1.366	-4.6	71	-0.06
12		1,3-Dichlorobenzene	1.481	1.591	-7.4	72	-0.06
13	C	1,4-Dichlorobenzene	1.524	1.603	-5.2	70	-0.06
14		1,2-Dichlorobenzene	1.419	1.512	-6.6	70	-0.06
15		Benzyl Alcohol	1.140	1.151	-1.0	67	-0.06
16		2,2'-oxybis(1-Chloropropane	1.998	2.005	-0.4	67	-0.06
17		2-Methylphenol	1.060	1.114	-5.1	70	-0.06
18		Hexachloroethane	0.525	0.570	-8.6	70	-0.06
19	P	n-Nitroso-di-n-propylamine	1.037	1.099	-6.0	67	-0.06
20		3+4-Methylphenols	1.413	1.550	-9.7	71	-0.06
21	I	Naphthalene-d8	1.000	1.000	0.0	63	-0.06
22		Acetophenone	0.452	0.472	-4.4	71	-0.06
23	S	Nitrobenzene-d5	0.348	0.386	-10.9	73	-0.06
24		Nitrobenzene	0.345	0.370	-7.2	73	-0.06
25		Isophorone	0.645	0.687	-6.5	70	-0.06
26	C	2-Nitrophenol	0.143	0.171	-19.6	75	-0.06
27		2,4-Dimethylphenol	0.286	0.296	-3.5	67	-0.06
28		bis(2-Chloroethoxy)methane	0.398	0.426	-7.0	69	-0.06
29	C	2,4-Dichlorophenol	0.254	0.277	-9.1	72	-0.07
30		1,2,4-Trichlorobenzene	0.279	0.303	-8.6	72	-0.06
31		Naphthalene	0.953	1.037	-8.8	73	-0.06
32		Benzoic acid	0.164	0.184	-12.2	68	-0.07
33		4-Chloroaniline	0.403	0.447	-10.9	75	-0.06
34	C	Hexachlorobutadiene	0.161	0.165	-2.5	69	-0.06
35		Caprolactam	0.082	0.089	-8.5	70	-0.08
36	C	4-Chloro-3-methylphenol	0.267	0.297	-11.2	68	-0.06
37		2-Methylnaphthalene	0.660	0.712	-7.9	71	-0.06
38	I	Acenaphthene-d10	1.000	1.000	0.0	62	-0.07
39		1,2,4,5-Tetrachlorobenzene	0.563	0.607	-7.8	72	-0.07
40	P	Hexachlorocyclopentadiene	0.283	0.146	48.4#	33#	-0.06
41	S	2,4,6-Tribromophenol	0.216	0.247	-14.4	71	-0.07
42	C	2,4,6-Trichlorophenol	0.387	0.416	-7.5	68	-0.07
43		2,4,5-Trichlorophenol	0.392	0.413	-5.4	67	-0.06
44	S	2-Fluorobiphenyl	1.436	1.592	-10.9	72	-0.06
45		1,1'-Biphenyl	1.590	1.715	-7.9	72	-0.06
46		2-Chloronaphthalene	1.240	1.357	-9.4	74	-0.06
47		2-Nitroaniline	0.359	0.405	-12.8	71	-0.06
48		Acenaphthylene	2.036	2.250	-10.5	73	-0.07

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49		Dimethylphthalate	1.468	1.558	-6.1	72	-0.07
50		2,6-Dinitrotoluene	0.290	0.339	-16.9	75	-0.06
51	C	Acenaphthene	1.214	1.288	-6.1	69	-0.06
52		3-Nitroaniline	0.362	0.408	-12.7	73	-0.07
53	P	2,4-Dinitrophenol	0.085	0.082	3.5	62	-0.06
54		Dibenzofuran	1.754	1.917	-9.3	71	-0.06
55	P	4-Nitrophenol	0.286	0.256	10.5	58	-0.06
56		2,4-Dinitrotoluene	0.383	0.422	-10.2	68	-0.07
57		Fluorene	1.440	1.601	-11.2	74	-0.06
58		2,3,4,6-Tetrachlorophenol	0.320	0.316	1.3	62	-0.06
59		Diethylphthalate	1.454	1.592	-9.5	72	-0.06
60		4-Chlorophenyl-phenylether	0.639	0.699	-9.4	72	-0.06
61		4-Nitroaniline	0.362	0.414	-14.4	72	-0.07
62		Azobenzene	1.433	1.519	-6.0	68	-0.06
63	I	Phenanthrene-d10	1.000	1.000	0.0	63	-0.07
64		4,6-Dinitro-2-methylphenol	0.080	0.089	-11.2	73	-0.07
65	c	n-Nitrosodiphenylamine	0.666	0.719	-8.0	70	-0.06
66		4-Bromophenyl-phenylether	0.208	0.220	-5.8	72	-0.06
67		Hexachlorobenzene	0.238	0.257	-8.0	74	-0.07
68		Atrazine	0.194	0.203	-4.6	71	-0.07
69	C	Pentachlorophenol	0.120	0.100	16.7	54	-0.07
70		Phenanthrene	1.119	1.172	-4.7	69	-0.06
71		Anthracene	1.098	1.162	-5.8	71	-0.07
72		Carbazole	1.046	1.132	-8.2	72	-0.06
73		Di-n-butylphthalate	1.255	1.361	-8.4	69	-0.07
74	C	Fluoranthene	1.166	1.251	-7.3	71	-0.07
75	I	Chrysene-d12	1.000	1.000	0.0	59	-0.09
76		Benzidine	0.539	0.653	-21.2	70	-0.07
77		Pyrene	1.152	1.268	-10.1	70	-0.07
78	S	Terphenyl-d14	0.835	0.917	-9.8	70	-0.07
79		Butylbenzylphthalate	0.504	0.584	-15.9	68	-0.07
80		Benzo(a)anthracene	1.095	1.175	-7.3	67	-0.09
81		3,3'-Dichlorobenzidine	0.390	0.439	-12.6	68	-0.09
82		Chrysene	1.053	1.142	-8.5	70	-0.09
83		Bis(2-ethylhexyl)phthalate	0.763	0.886	-16.1	68	-0.08
84	c	Di-n-octyl phthalate	1.344	1.499	-11.5	69	-0.13
85		Indeno(1,2,3-cd)pyrene	1.342	1.487	-10.8	69	-0.24
86	I	Perylene-d12	1.000	1.000	0.0	63	-0.18
87		Benzo(b)fluoranthene	1.095	1.142	-4.3	68	-0.15
88		Benzo(k)fluoranthene	1.060	1.184	-11.7	72	-0.15
89	C	Benzo(a)pyrene	1.008	1.128	-11.9	73	-0.17
90		Dibenzo(a,h)anthracene	1.071	1.202	-12.2	73	-0.25
91		Benzo(g,h,i)perylene	1.074	1.202	-11.9	73	-0.25

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