Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF060315\
Data File : BF079674.D
Acq On : 3 Jun 2015 21:07
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 05 11:39:26 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 05 10:52:02 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	95	0.00
2		1,4-Dioxane	0.510	0.508	0.4	95	0.00
3		Pyridine	1.466	1.459	0.5	104	0.00
4		n-Nitrosodimethylamine	0.654	0.691	-5.7	99	0.00
5	S	2-Fluorophenol	1.256	1.111	11.5	94	0.00
6		Aniline	2.079	2.128	-2.4	94	0.00
7	S	Phenol-d6	1.679	1.558	7.2	96	0.00
8		2-Chlorophenol	1.282	1.348	-5.1	98	0.00
9		Benzaldehyde	0.840	0.872	-3.8	96	0.00
10	C	Phenol	1.553	1.645	-5.9	97	0.00
11		bis(2-Chloroethyl)ether	1.238	1.275	-3.0	94	0.00
12		1,3-Dichlorobenzene	1.466	1.498	-2.2	98	0.00
13	C	1,4-Dichlorobenzene	1.499	1.516	-1.1	94	0.00
14		1,2-Dichlorobenzene	1.357	1.388	-2.3	96	-0.01
15		Benzyl Alcohol	1.056	1.063	-0.7	96	0.00
16		2,2'-oxybis(1-Chloropropane	2.065	2.032	1.6	91	0.00
17		2-Methylphenol	1.059	1.074	-1.4	96	0.00
18		Hexachloroethane	0.491	0.490	0.2	91	0.00
19	P	n-Nitroso-di-n-propylamine	0.991	1.051	-6.1	90	0.00
20		3+4-Methylphenols	1.402	1.431	-2.1	96	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
22		Acetophenone	0.468	0.475	-1.5	94	0.00
23	S	Nitrobenzene-d5	0.390	0.346	11.3	90	0.00
24		Nitrobenzene	0.335	0.348	-3.9	97	0.00
25		Isophorone	0.682	0.709	-4.0	94	0.00
26	C	2-Nitrophenol	0.111	0.112	-0.9	91	0.00
27		2,4-Dimethylphenol	0.312	0.284	9.0	87	0.00
28		bis(2-Chloroethoxy)methane	0.392	0.393	-0.3	97	0.00
29	C	2,4-Dichlorophenol	0.278	0.297	-6.8	99	0.00
30		1,2,4-Trichlorobenzene	0.314	0.311	1.0	91	0.00
31		Naphthalene	1.046	1.040	0.6	98	0.00
32		Benzoic acid	0.086	0.099	-15.1	105	0.00
33		4-Chloroaniline	0.569	0.585	-2.8	97	0.00
34	C	Hexachlorobutadiene	0.176	0.172	2.3	92	-0.01
35		Caprolactam	0.077	0.085	-10.4	97	-0.01
36	C	4-Chloro-3-methylphenol	0.284	0.301	-6.0	102	0.00
37		2-Methylnaphthalene	0.698	0.725	-3.9	99	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
39		1,2,4,5-Tetrachlorobenzene	0.642	0.664	-3.4	94	0.00
40	P	Hexachlorocyclopentadiene	0.306	0.276	9.8	80	0.00
41	S	2,4,6-Tribromophenol	0.165	0.170	-3.0	90	-0.01
42	C	2,4,6-Trichlorophenol	0.349	0.373	-6.9	91	0.00
43		2,4,5-Trichlorophenol	0.387	0.445	-15.0	98	0.00
44	S	2-Fluorobiphenyl	1.665	1.438	13.6	96	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.542	1.582	-2.6	94	-0.01
46		2-Chloronaphthalene	1.208	1.264	-4.6	98	-0.01
47		2-Nitroaniline	0.249	0.301	-20.9	104	0.00
48		Acenaphthylene	2.003	2.081	-3.9	94	0.00
49		Dimethylphthalate	1.465	1.520	-3.8	100	0.00
50		2,6-Dinitrotoluene	0.233	0.272	-16.7	98	0.00
51	C	Acenaphthene	1.214	1.262	-4.0	94	0.00
52		3-Nitroaniline	0.287	0.356	-24.0	106	0.00
53	P	2,4-Dinitrophenol	0.066	0.067	-1.5	92	0.00
54		Dibenzofuran	1.762	1.864	-5.8	96	0.00
55	P	4-Nitrophenol	0.232	0.245	-5.6	93	0.00
56		2,4-Dinitrotoluene	0.280	0.311	-11.1	93	0.00
57		Fluorene	1.440	1.561	-8.4	102	-0.01
58		2,3,4,6-Tetrachlorophenol	0.268	0.318	-18.7	100	-0.01
59		Diethylphthalate	1.407	1.595	-13.4	107	0.00
60		4-Chlorophenyl-phenylether	0.650	0.698	-7.4	94	0.00
61		4-Nitroaniline	0.283	0.351	-24.0	107	0.00
62		Azobenzene	1.316	1.382	-5.0	90	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	-0.02
64		4,6-Dinitro-2-methylphenol	0.052	0.055	-5.8	99	0.00
65	c	n-Nitrosodiphenylamine	0.628	0.625	0.5	95	-0.01
66		4-Bromophenyl-phenylether	0.201	0.207	-3.0	102	-0.01
67		Hexachlorobenzene	0.236	0.233	1.3	97	-0.01
68		Atrazine	0.161	0.182	-13.0	95	-0.01
69	C	Pentachlorophenol	0.086	0.104	-20.9#	110	-0.02
70		Phenanthrene	1.106	1.153	-4.2	100	-0.02
71		Anthracene	1.113	1.133	-1.8	97	-0.02
72		Carbazole	1.022	1.056	-3.3	102	-0.02
73		Di-n-butylphthalate	1.120	1.203	-7.4	98	-0.05
74	C	Fluoranthene	1.192	1.228	-3.0	102	-0.07
75	I	Chrysene-d12	1.000	1.000	0.0	101	-0.19
76		Benzidine	0.455	0.399	12.3	77	-0.08
77		Pyrene	1.152	1.180	-2.4	98	-0.09
78	S	Terphenyl-d14	0.948	0.849	10.4	100	-0.10
79		Butylbenzylphthalate	0.357	0.441	-23.5	106	-0.15
80		Benzo(a)anthracene	1.094	1.150	-5.1	100	-0.19
81		3,3'-Dichlorobenzidine	0.327	0.388	-18.7	104	-0.19
82		Chrysene	1.100	1.131	-2.8	100	-0.19
83		Bis(2-ethylhexyl)phthalate	0.568	0.708	-24.6	106	-0.22
84	c	Di-n-octyl phthalate	0.844	1.095	-29.7#	108	-0.29
85		Indeno(1,2,3-cd)pyrene	1.076	1.219	-13.3	107	-0.32
86	I	Perylene-d12	1.000	1.000	0.0	102	-0.30
87		Benzo(b)fluoranthene	1.034	1.167	-12.9	103	-0.29

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88	Benzo(k)fluoranthene	1.211	1.248	-3.1	100	-0.29
89 C	Benzo(a)pyrene	1.011	1.133	-12.1	103	-0.29
90	Dibenzo(a,h)anthracene	0.947	1.107	-16.9	108	-0.32
91	Benzo(g,h,i)perylene	1.004	1.118	-11.4	107	-0.33

(#) = Out of Range SPCC's out = 0 CCC's out = 2