Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061715\
Data File : BF080031.D
Acq On : 17 Jun 2015 18:37
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 18 00:54:19 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 17 16:41:53 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	108	0.00
2		1,4-Dioxane	0.537	0.522	2.8	105	0.00
3		Pyridine	1.428	1.429	-0.1	108	0.01
4		n-Nitrosodimethylamine	0.679	0.661	2.7	99	0.00
5	S	2-Fluorophenol	1.130	1.180	-4.4	110	0.00
6		Aniline	2.068	2.182	-5.5	109	0.00
7	S	Phenol-d6	1.503	1.566	-4.2	105	0.00
8		2-Chlorophenol	1.306	1.350	-3.4	107	0.00
9		Benzaldehyde	0.868	0.826	4.8	99	0.00
10	C	Phenol	1.641	1.704	-3.8	107	-0.01
11		bis(2-Chloroethyl)ether	1.302	1.364	-4.8	109	0.00
12		1,3-Dichlorobenzene	1.569	1.604	-2.2	106	0.00
13	C	1,4-Dichlorobenzene	1.492	1.538	-3.1	108	0.00
14		1,2-Dichlorobenzene	1.452	1.473	-1.4	106	0.00
15		Benzyl Alcohol	1.229	1.265	-2.9	106	0.00
16		2,2'-oxybis(1-Chloropropane	1.932	1.898	1.8	107	0.00
17		2-Methylphenol	1.115	1.170	-4.9	107	0.00
18		Hexachloroethane	0.497	0.520	-4.6	107	0.00
19	P	n-Nitroso-di-n-propylamine	1.044	1.008	3.4	108	-0.01
20		3+4-Methylphenols	1.440	1.467	-1.9	105	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
22		Acetophenone	0.466	0.491	-5.4	108	0.00
23	S	Nitrobenzene-d5	0.344	0.352	-2.3	108	0.00
24		Nitrobenzene	0.355	0.353	0.6	106	0.00
25		Isophorone	0.691	0.703	-1.7	108	0.00
26	C	2-Nitrophenol	0.148	0.148	0.0	106	-0.01
27		2,4-Dimethylphenol	0.309	0.303	1.9	109	0.00
28		bis(2-Chloroethoxy)methane	0.406	0.421	-3.7	110	0.00
29	C	2,4-Dichlorophenol	0.265	0.265	0.0	105	0.00
30		1,2,4-Trichlorobenzene	0.301	0.309	-2.7	111	0.00
31		Naphthalene	1.046	1.051	-0.5	106	0.00
32		Benzoic acid	0.187	0.180	3.7	102	-0.01
33		4-Chloroaniline	0.448	0.437	2.5	109	0.00
34	C	Hexachlorobutadiene	0.185	0.184	0.5	112	0.00
35		Caprolactam	0.086	0.091	-5.8	107	-0.01
36	C	4-Chloro-3-methylphenol	0.299	0.311	-4.0	107	0.00
37		2-Methylnaphthalene	0.676	0.684	-1.2	107	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
39		1,2,4,5-Tetrachlorobenzene	0.582	0.597	-2.6	109	-0.01
40	P	Hexachlorocyclopentadiene	0.260	0.279	-7.3	112	0.00
41	S	2,4,6-Tribromophenol	0.196	0.199	-1.5	108	-0.01
42	C	2,4,6-Trichlorophenol	0.396	0.405	-2.3	107	0.00
43		2,4,5-Trichlorophenol	0.456	0.474	-3.9	108	0.00
44	S	2-Fluorobiphenyl	1.402	1.389	0.9	106	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.627	1.575	3.2	105	0.00
46		2-Chloronaphthalene	1.320	1.359	-3.0	108	0.00
47		2-Nitroaniline	0.362	0.388	-7.2	107	0.00
48		Acenaphthylene	2.204	2.322	-5.4	108	0.00
49		Dimethylphthalate	1.504	1.487	1.1	108	0.00
50		2,6-Dinitrotoluene	0.301	0.337	-12.0	111	0.00
51	C	Acenaphthene	1.348	1.374	-1.9	107	0.00
52		3-Nitroaniline	0.348	0.383	-10.1	112	0.00
53	P	2,4-Dinitrophenol	0.106	0.117	-10.4	113	0.00
54		Dibenzofuran	1.913	1.915	-0.1	106	0.00
55	P	4-Nitrophenol	0.278	0.274	1.4	108	-0.01
56		2,4-Dinitrotoluene	0.383	0.394	-2.9	108	0.00
57		Fluorene	1.518	1.486	2.1	105	0.00
58		2,3,4,6-Tetrachlorophenol	0.385	0.409	-6.2	109	0.00
59		Diethylphthalate	1.514	1.596	-5.4	107	0.00
60		4-Chlorophenyl-phenylether	0.729	0.746	-2.3	111	0.00
61		4-Nitroaniline	0.359	0.397	-10.6	113	0.00
62		Azobenzene	1.541	1.619	-5.1	107	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	108	0.00
64		4,6-Dinitro-2-methylphenol	0.094	0.107	-13.8	113	0.00
65	c	n-Nitrosodiphenylamine	0.651	0.717	-10.1	111	0.00
66		4-Bromophenyl-phenylether	0.213	0.222	-4.2	108	0.00
67		Hexachlorobenzene	0.236	0.243	-3.0	106	0.01
68		Atrazine	0.206	0.212	-2.9	103	0.01
69	C	Pentachlorophenol	0.134	0.146	-9.0	118	0.00
70		Phenanthrene	1.211	1.221	-0.8	107	0.01
71		Anthracene	1.166	1.227	-5.2	112	0.00
72		Carbazole	1.113	1.151	-3.4	107	0.01
73		Di-n-butylphthalate	1.286	1.347	-4.7	115	0.02
74	C	Fluoranthene	1.290	1.357	-5.2	114	0.05
75	I	Chrysene-d12	1.000	1.000	0.0	108	0.14
76		Benzidine	0.559	0.469	16.1	88	0.06
77		Pyrene	1.231	1.219	1.0	106	0.06
78	S	Terphenyl-d14	0.856	0.874	-2.1	111	0.07
79		Butylbenzylphthalate	0.495	0.510	-3.0	106	0.09
80		Benzo(a)anthracene	1.218	1.236	-1.5	110	0.14
81		3,3'-Dichlorobenzidine	0.420	0.438	-4.3	111	0.14
82		Chrysene	1.124	1.146	-2.0	109	0.14
83		Bis(2-ethylhexyl)phthalate	0.782	0.796	-1.8	106	0.14
84	c	Di-n-octyl phthalate	1.339	1.340	-0.1	105	0.19
85		Indeno(1,2,3-cd)pyrene	1.417	1.414	0.2	108	0.22
86	I	Perylene-d12	1.000	1.000	0.0	110	0.21
87		Benzo(b)fluoranthene	1.211	1.244	-2.7	109	0.19

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88	Benzo(k)fluoranthene	1.233	1.183	4.1	106	0.19
89 C	Benzo(a)pyrene	1.150	1.150	0.0	109	0.21
90	Dibenzo(a,h)anthracene	1.177	1.174	0.3	109	0.22
91	Benzo(g,h,i)perylene	1.188	1.184	0.3	108	0.23

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