Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF072216\
Data File : BF089187.D
Acq On : 22 Jul 2016 10:11
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 22 11:11:51 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 20 19:03:15 2016
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	99	-0.01
2		1,4-Dioxane	0.571	0.587	-2.8	102	-0.02
3		Pyridine	1.450	1.425	1.7	99	-0.02
4		n-Nitrosodimethylamine	0.611	0.586	4.1	94	-0.02
5	S	2-Fluorophenol	1.315	1.298	1.3	99	-0.01
6		Aniline	2.291	2.137	6.7	96	-0.01
7	S	Phenol-d6	1.696	1.641	3.2	99	-0.02
8		2-Chlorophenol	1.459	1.487	-1.9	105	-0.02
9		Benzaldehyde	0.906	0.946	-4.4	106	-0.01
10	C	Phenol	1.954	1.929	1.3	97	-0.01
11		bis(2-Chloroethyl)ether	1.467	1.581	-7.8	106	-0.01
12		1,3-Dichlorobenzene	1.611	1.542	4.3	102	-0.01
13	C	1,4-Dichlorobenzene	1.623	1.602	1.3	104	-0.01
14		1,2-Dichlorobenzene	1.530	1.621	-5.9	104	-0.01
15		Benzyl Alcohol	1.242	1.224	1.4	95	-0.01
16		2,2'-oxybis(1-Chloropropane	1.672	1.733	-3.6	105	-0.01
17		2-Methylphenol	1.149	1.120	2.5	98	-0.01
18		Hexachloroethane	0.610	0.606	0.7	104	-0.01
19	P	n-Nitroso-di-n-propylamine	1.109	1.160	-4.6	101	-0.01
20		3+4-Methylphenols	1.560	1.539	1.3	97	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	100	-0.01
22		Acetophenone	0.534	0.543	-1.7	101	-0.01
23	S	Nitrobenzene-d5	0.376	0.374	0.5	104	-0.01
24		Nitrobenzene	0.397	0.418	-5.3	103	-0.01
25		Isophorone	0.694	0.682	1.7	101	-0.01
26	C	2-Nitrophenol	0.185	0.182	1.6	104	0.00
27		2,4-Dimethylphenol	0.312	0.297	4.8	102	-0.01
28		bis(2-Chloroethoxy)methane	0.434	0.457	-5.3	100	-0.01
29	C	2,4-Dichlorophenol	0.281	0.291	-3.6	105	-0.01
30		1,2,4-Trichlorobenzene	0.311	0.317	-1.9	103	-0.01
31		Naphthalene	1.079	1.007	6.7	100	-0.01
32		Benzoic acid	0.240	0.243	-1.3	98	-0.01
33		4-Chloroaniline	0.454	0.458	-0.9	103	-0.01
34	C	Hexachlorobutadiene	0.173	0.178	-2.9	103	-0.01
35		Caprolactam	0.087	0.086	1.1	100	-0.01
36	C	4-Chloro-3-methylphenol	0.339	0.357	-5.3	109	-0.01
37		2-Methylnaphthalene	0.688	0.699	-1.6	101	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	102	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.743	0.700	5.8	103	-0.01
40	P	Hexachlorocyclopentadiene	0.201	0.272	-35.3#	132	-0.01
41	S	2,4,6-Tribromophenol	0.180	0.178	1.1	105	-0.01
42	C	2,4,6-Trichlorophenol	0.416	0.394	5.3	94	-0.01
43		2,4,5-Trichlorophenol	0.418	0.423	-1.2	101	-0.01
44	S	2-Fluorobiphenyl	1.452	1.538	-5.9	103	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.801	1.648	8.5	102	-0.01
46		2-Chloronaphthalene	1.369	1.326	3.1	104	-0.01
47		2-Nitroaniline	0.470	0.446	5.1	103	-0.01
48		Acenaphthylene	2.152	2.154	-0.1	107	-0.01
49		Dimethylphthalate	1.535	1.477	3.8	95	-0.01
50		2,6-Dinitrotoluene	0.346	0.344	0.6	95	-0.01
51	C	Acenaphthene	1.273	1.283	-0.8	107	-0.01
52		3-Nitroaniline	0.411	0.391	4.9	91	-0.01
53	P	2,4-Dinitrophenol	0.141	0.167	-18.4	111	-0.01
54		Dibenzofuran	1.720	1.686	2.0	102	-0.01
55	P	4-Nitrophenol	0.249	0.231	7.2	87	-0.01
56		2,4-Dinitrotoluene	0.446	0.436	2.2	100	-0.01
57		Fluorene	1.460	1.508	-3.3	97	-0.01
58		2,3,4,6-Tetrachlorophenol	0.300	0.331	-10.3	107	-0.01
59		Diethylphthalate	1.522	1.441	5.3	101	-0.01
60		4-Chlorophenyl-phenylether	0.667	0.624	6.4	100	-0.01
61		4-Nitroaniline	0.400	0.396	1.0	103	-0.01
62		Azobenzene	1.654	1.581	4.4	105	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	98	-0.01
64		4,6-Dinitro-2-methylphenol	0.116	0.123	-6.0	97	-0.01
65	c	n-Nitrosodiphenylamine	0.668	0.658	1.5	103	-0.01
66		4-Bromophenyl-phenylether	0.198	0.206	-4.0	102	-0.01
67		Hexachlorobenzene	0.213	0.203	4.7	99	-0.02
68		Atrazine	0.209	0.215	-2.9	101	-0.01
69	C	Pentachlorophenol	0.098	0.131	-33.7#	125	-0.01
70		Phenanthrene	1.097	1.096	0.1	106	-0.01
71		Anthracene	1.140	1.168	-2.5	99	-0.01
72		Carbazole	1.002	1.029	-2.7	96	-0.01
73		Di-n-butylphthalate	1.239	1.142	7.8	99	-0.01
74	C	Fluoranthene	1.153	1.103	4.3	103	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	95	-0.01
76		Benzidine	0.688	1.247	-81.3#	173#	-0.01
77		Pyrene	1.652	1.767	-7.0	100	-0.01
78	S	Terphenyl-d14	0.922	0.889	3.6	102	-0.01
79		Butylbenzylphthalate	0.699	0.732	-4.7	98	-0.01
80		Benzo(a)anthracene	1.274	1.267	0.5	97	-0.01
81		3,3'-Dichlorobenzidine	0.423	0.472	-11.6	95	-0.01
82		Chrysene	1.217	1.309	-7.6	98	-0.01
83		Bis(2-ethylhexyl)phthalate	0.927	0.984	-6.1	105	-0.01
84	c	Di-n-octyl phthalate	1.529	1.645	-7.6	104	0.00
85		Indeno(1,2,3-cd)pyrene	0.884	0.858	2.9	92	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	92	-0.01
87		Benzo(b)fluoranthene	1.449	1.568	-8.2	88	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.036	0.961	7.2	103	-0.01
89 C	Benzo(a)pyrene	1.125	1.118	0.6	95	-0.01
90	Dibenzo(a,h)anthracene	0.888	0.875	1.5	93	-0.02
91	Benzo(g,h,i)perylene	0.895	0.866	3.2	92	-0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 1