Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102715\
Data File : BF082551.D
Acq On : 27 Oct 2015 17:19
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 27 17:49:33 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 27 02:43:58 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	93	-0.02
2		1,4-Dioxane	0.508	0.454	10.6	90	-0.03
3		Pyridine	1.185	1.173	1.0	88	-0.03
4		n-Nitrosodimethylamine	0.521	0.527	-1.2	93	-0.03
5	S	2-Fluorophenol	1.013	1.053	-3.9	98	-0.01
6		Aniline	1.664	1.641	1.4	94	-0.01
7	S	Phenol-d6	1.320	1.289	2.3	96	0.00
8		2-Chlorophenol	1.188	1.241	-4.5	98	-0.01
9		Benzaldehyde	0.688	0.690	-0.3	95	-0.01
10	C	Phenol	1.499	1.532	-2.2	106	-0.01
11		bis(2-Chloroethyl)ether	1.190	1.210	-1.7	102	-0.01
12		1,3-Dichlorobenzene	1.417	1.411	0.4	98	0.07
13	C	1,4-Dichlorobenzene	1.458	1.430	1.9	95	-0.01
14		1,2-Dichlorobenzene	1.354	1.438	-6.2	111	-0.17
15		Benzyl Alcohol	0.919	0.951	-3.5	100	-0.01
16		2,2'-oxybis(1-Chloropropane	1.611	1.551	3.7	92	-0.02
17		2-Methylphenol	0.953	0.956	-0.3	92	-0.01
18		Hexachloroethane	0.496	0.487	1.8	93	-0.02
19	P	n-Nitroso-di-n-propylamine	0.860	0.900	-4.7	106	-0.01
20		3+4-Methylphenols	1.218	1.238	-1.6	97	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	97	-0.02
22		Acetophenone	0.434	0.423	2.5	96	-0.02
23	S	Nitrobenzene-d5	0.310	0.306	1.3	99	-0.01
24		Nitrobenzene	0.346	0.342	1.2	93	-0.01
25		Isophorone	0.616	0.609	1.1	99	-0.01
26	C	2-Nitrophenol	0.167	0.170	-1.8	100	-0.02
27		2,4-Dimethylphenol	0.272	0.272	0.0	101	-0.01
28		bis(2-Chloroethoxy)methane	0.361	0.346	4.2	98	-0.02
29	C	2,4-Dichlorophenol	0.271	0.285	-5.2	102	-0.01
30		1,2,4-Trichlorobenzene	0.307	0.294	4.2	97	-0.02
31		Naphthalene	0.914	0.880	3.7	100	-0.02
32		Benzoic acid	0.152	0.130	14.5	80	0.00
33		4-Chloroaniline	0.354	0.350	1.1	92	-0.01
34	C	Hexachlorobutadiene	0.177	0.182	-2.8	106	-0.01
35		Caprolactam	0.075	0.073	2.7	99	-0.01
36	C	4-Chloro-3-methylphenol	0.264	0.274	-3.8	97	-0.01
37		2-Methylnaphthalene	0.617	0.631	-2.3	102	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	99	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.604	0.601	0.5	100	-0.02
40	P	Hexachlorocyclopentadiene	0.115	0.070	39.1#	55	-0.02
41	S	2,4,6-Tribromophenol	0.150	0.143	4.7	93	-0.02
42	C	2,4,6-Trichlorophenol	0.376	0.384	-2.1	99	-0.01
43		2,4,5-Trichlorophenol	0.395	0.394	0.3	94	-0.01
44	S	2-Fluorobiphenyl	1.258	1.305	-3.7	110	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.488	1.480	0.5	92	-0.01
46		2-Chloronaphthalene	1.173	1.203	-2.6	100	-0.01
47		2-Nitroaniline	0.341	0.345	-1.2	103	-0.01
48		Acenaphthylene	1.879	1.914	-1.9	106	-0.01
49		Dimethylphthalate	1.392	1.435	-3.1	101	-0.01
50		2,6-Dinitrotoluene	0.305	0.311	-2.0	94	-0.01
51	C	Acenaphthene	1.078	1.107	-2.7	106	-0.01
52		3-Nitroaniline	0.331	0.355	-7.3	98	-0.01
53	P	2,4-Dinitrophenol	0.112	0.102	8.9	88	-0.01
54		Dibenzofuran	1.572	1.567	0.3	104	-0.01
55	P	4-Nitrophenol	0.113	0.082	27.4#	65	0.00
56		2,4-Dinitrotoluene	0.370	0.392	-5.9	108	-0.01
57		Fluorene	1.291	1.353	-4.8	111	-0.01
58		2,3,4,6-Tetrachlorophenol	0.302	0.296	2.0	96	-0.01
59		Diethylphthalate	1.354	1.281	5.4	93	-0.02
60		4-Chlorophenyl-phenylether	0.607	0.627	-3.3	108	-0.02
61		4-Nitroaniline	0.332	0.324	2.4	88	-0.01
62		Azobenzene	1.308	1.310	-0.2	106	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	-0.01
64		4,6-Dinitro-2-methylphenol	0.117	0.124	-6.0	101	-0.01
65	c	n-Nitrosodiphenylamine	0.605	0.615	-1.7	106	-0.01
66		4-Bromophenyl-phenylether	0.188	0.200	-6.4	109	-0.01
67		Hexachlorobenzene	0.200	0.197	1.5	93	-0.02
68		Atrazine	0.187	0.198	-5.9	101	-0.01
69	C	Pentachlorophenol	0.068	0.057	16.2	74	-0.01
70		Phenanthrene	1.029	1.054	-2.4	111	-0.01
71		Anthracene	1.037	1.047	-1.0	96	-0.02
72		Carbazole	0.914	0.945	-3.4	100	-0.01
73		Di-n-butylphthalate	1.090	1.144	-5.0	110	-0.01
74	C	Fluoranthene	1.057	1.090	-3.1	107	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	89	-0.02
76		Benzidine	0.549	0.496	9.7	79	-0.02
77		Pyrene	1.230	1.379	-12.1	106	-0.01
78	S	Terphenyl-d14	0.694	0.718	-3.5	92	-0.02
79		Butylbenzylphthalate	0.517	0.576	-11.4	105	-0.01
80		Benzo(a)anthracene	1.072	1.033	3.6	91	-0.02
81		3,3'-Dichlorobenzidine	0.372	0.376	-1.1	87	-0.02
82		Chrysene	1.013	0.980	3.3	91	-0.02
83		Bis(2-ethylhexyl)phthalate	0.693	0.680	1.9	88	-0.02
84	c	Di-n-octyl phthalate	1.143	0.947	17.1	75	-0.02
85		Indeno(1,2,3-cd)pyrene	0.931	0.751	19.3	76	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	68	-0.02
87		Benzo(b)fluoranthene	1.188	1.100	7.4	66	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	0.968	1.182	-22.1	86	-0.03
89 C	Benzo(a)pyrene	1.029	1.037	-0.8	71	-0.03
90	Dibenzo(a,h)anthracene	0.861	0.925	-7.4	76	-0.05
91	Benzo(g,h,i)perylene	0.852	0.921	-8.1	79	-0.05

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