Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090916\
Data File : BF090049.D
Acq On : 9 Sep 2016 16:03
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 09 17:57:51 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 08 16:07:52 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	94	-0.01
2		1,4-Dioxane	0.592	0.555	6.2	94	-0.02
3		Pyridine	1.579	1.457	7.7	83	-0.02
4		n-Nitrosodimethylamine	0.779	0.673	13.6	79	-0.02
5	S	2-Fluorophenol	1.250	1.250	0.0	90	-0.01
6		Aniline	2.072	1.862	10.1	89	-0.01
7	S	Phenol-d6	1.518	1.423	6.3	87	-0.01
8		2-Chlorophenol	1.356	1.347	0.7	92	-0.01
9		Benzaldehyde	0.980	0.868	11.4	83	-0.01
10	C	Phenol	1.766	1.567	11.3	86	-0.01
11		bis(2-Chloroethyl)ether	1.336	1.272	4.8	95	-0.01
12		1,3-Dichlorobenzene	1.514	1.479	2.3	94	-0.01
13	C	1,4-Dichlorobenzene	1.549	1.548	0.1	98	-0.01
14		1,2-Dichlorobenzene	1.389	1.411	-1.6	93	-0.01
15		Benzyl Alcohol	0.951	0.914	3.9	90	-0.01
16		2,2'-oxybis(1-Chloropropane	2.547	2.452	3.7	88	-0.01
17		2-Methylphenol	1.083	1.096	-1.2	95	-0.01
18		Hexachloroethane	0.531	0.525	1.1	93	-0.02
19	P	n-Nitroso-di-n-propylamine	0.989	0.923	6.7	83	-0.02
20		3+4-Methylphenols	1.312	1.342	-2.3	108	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	96	-0.02
22		Acetophenone	0.449	0.433	3.6	96	-0.02
23	S	Nitrobenzene-d5	0.345	0.341	1.2	96	-0.01
24		Nitrobenzene	0.367	0.323	12.0	91	-0.02
25		Isophorone	0.701	0.665	5.1	94	-0.01
26	C	2-Nitrophenol	0.175	0.188	-7.4	104	-0.01
27		2,4-Dimethylphenol	0.324	0.314	3.1	103	-0.01
28		bis(2-Chloroethoxy)methane	0.423	0.413	2.4	91	-0.01
29	C	2,4-Dichlorophenol	0.291	0.286	1.7	100	-0.01
30		1,2,4-Trichlorobenzene	0.318	0.308	3.1	91	-0.01
31		Naphthalene	0.996	0.883	11.3	93	-0.01
32		Benzoic acid	0.218	0.229	-5.0	102	-0.01
33		4-Chloroaniline	0.402	0.298	25.9#	70	-0.01
34	C	Hexachlorobutadiene	0.186	0.176	5.4	94	-0.01
35		Caprolactam	0.092	0.088	4.3	88	-0.01
36	C	4-Chloro-3-methylphenol	0.308	0.289	6.2	91	-0.01
37		2-Methylnaphthalene	0.658	0.647	1.7	93	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	93	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.586	0.556	5.1	100	-0.01
40	P	Hexachlorocyclopentadiene	0.300	0.278	7.3	84	-0.02
41	S	2,4,6-Tribromophenol	0.197	0.197	0.0	92	-0.01
42	C	2,4,6-Trichlorophenol	0.406	0.408	-0.5	88	-0.01
43		2,4,5-Trichlorophenol	0.390	0.431	-10.5	111	-0.01
44	S	2-Fluorobiphenyl	1.218	1.150	5.6	91	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.610	1.636	-1.6	105	-0.01
46		2-Chloronaphthalene	1.139	1.111	2.5	94	-0.01
47		2-Nitroaniline	0.372	0.367	1.3	95	-0.01
48		Acenaphthylene	1.884	1.779	5.6	90	-0.01
49		Dimethylphthalate	1.446	1.458	-0.8	107	-0.01
50		2,6-Dinitrotoluene	0.322	0.357	-10.9	106	-0.01
51	C	Acenaphthene	1.193	1.140	4.4	93	-0.01
52		3-Nitroaniline	0.367	0.321	12.5	91	-0.01
53	P	2,4-Dinitrophenol	0.130	0.159	-22.3	103	-0.01
54		Dibenzofuran	1.599	1.458	8.8	87	-0.01
55	P	4-Nitrophenol	0.277	0.262	5.4	81	-0.01
56		2,4-Dinitrotoluene	0.408	0.416	-2.0	85	0.00
57		Fluorene	1.169	1.190	-1.8	91	-0.01
58		2,3,4,6-Tetrachlorophenol	0.334	0.365	-9.3	106	-0.01
59		Diethylphthalate	1.360	1.343	1.3	98	0.00
60		4-Chlorophenyl-phenylether	0.550	0.556	-1.1	93	-0.01
61		4-Nitroaniline	0.360	0.369	-2.5	88	-0.01
62		Azobenzene	1.358	1.286	5.3	95	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	94	-0.01
64		4,6-Dinitro-2-methylphenol	0.128	0.125	2.3	100	-0.01
65	c	n-Nitrosodiphenylamine	0.601	0.590	1.8	98	0.00
66		4-Bromophenyl-phenylether	0.201	0.197	2.0	95	-0.01
67		Hexachlorobenzene	0.230	0.231	-0.4	106	0.00
68		Atrazine	0.201	0.180	10.4	91	-0.01
69	C	Pentachlorophenol	0.124	0.139	-12.1	94	0.00
70		Phenanthrene	0.996	0.900	9.6	86	-0.01
71		Anthracene	1.010	1.042	-3.2	110	0.00
72		Carbazole	0.947	0.971	-2.5	101	0.00
73		Di-n-butylphthalate	1.104	1.069	3.2	97	-0.01
74	C	Fluoranthene	1.044	0.978	6.3	86	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	75	-0.01
76		Benzidine	0.758	0.412	45.6#	41#	-0.01
77		Pyrene	1.410	1.665	-18.1	97	-0.01
78	S	Terphenyl-d14	0.810	0.985	-21.6	108	0.00
79		Butylbenzylphthalate	0.569	0.665	-16.9	90	0.00
80		Benzo(a)anthracene	1.225	1.198	2.2	75	-0.01
81		3,3'-Dichlorobenzidine	0.442	0.363	17.9	63	-0.01
82		Chrysene	1.144	0.989	13.5	68	-0.01
83		Bis(2-ethylhexyl)phthalate	0.787	0.853	-8.4	81	-0.01
84	c	Di-n-octyl phthalate	1.270	1.217	4.2	67	-0.01
85		Indeno(1,2,3-cd)pyrene	0.846	1.136	-34.3#	101	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	74	-0.01
87		Benzo(b)fluoranthene	1.253	1.213	3.2	63	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.063	0.868	18.3	71	-0.01
89 C	Benzo(a)pyrene	1.073	1.020	4.9	69	-0.01
90	Dibenzo(a,h)anthracene	0.787	1.081	-37.4#	101	-0.01
91	Benzo(g,h,i)perylene	0.792	1.112	-40.4#	105	-0.01

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