Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF101818\
Data File : BF109965.D
Acq On : 18 Oct 2018 18:10
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 19 03:59:56 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 16 14:58:24 2018
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	86	0.00
2		1,4-Dioxane	0.717	0.651	9.2	78	0.00
3		Pyridine	1.598	1.485	7.1	79	0.00
4		n-Nitrosodimethylamine	0.652	0.611	6.3	79	0.00
5	S	2-Fluorophenol	1.262	1.218	3.5	84	0.00
6		Aniline	2.111	2.034	3.6	82	0.00
7	S	Phenol-d6	1.567	1.528	2.5	85	-0.01
8		2-Chlorophenol	1.413	1.402	0.8	85	0.00
9		Benzaldehyde	1.018	0.985	3.2	84	0.00
10	C	Phenol	1.692	1.637	3.3	84	0.00
11		bis(2-Chloroethyl)ether	1.414	1.342	5.1	82	0.00
12		1,3-Dichlorobenzene	1.550	1.512	2.5	84	0.00
13	C	1,4-Dichlorobenzene	1.561	1.528	2.1	84	0.00
14		1,2-Dichlorobenzene	1.437	1.417	1.4	85	0.00
15		Benzyl Alcohol	1.205	1.216	-0.9	86	0.00
16		2,2'-oxybis(1-Chloropropane	2.321	2.167	6.6	80	0.00
17		2-Methylphenol	1.140	1.120	1.8	85	0.00
18		Hexachloroethane	0.552	0.563	-2.0	87	0.00
19	P	n-Nitroso-di-n-propylamine	1.006	0.962	4.4	84	0.00
20		3+4-Methylphenols	1.446	1.430	1.1	84	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
22		Acetophenone	0.464	0.434	6.5	83	0.00
23	S	Nitrobenzene-d5	0.297	0.327	-10.1	92	-0.01
24		Nitrobenzene	0.323	0.342	-5.9	89	0.00
25		Isophorone	0.583	0.555	4.8	84	0.00
26	C	2-Nitrophenol	0.134	0.177	-32.1#	112	0.00
27		2,4-Dimethylphenol	0.281	0.270	3.9	84	-0.01
28		bis(2-Chloroethoxy)methane	0.400	0.373	6.8	82	0.00
29	C	2,4-Dichlorophenol	0.273	0.275	-0.7	88	0.00
30		1,2,4-Trichlorobenzene	0.305	0.303	0.7	86	0.00
31		Naphthalene	0.917	0.889	3.1	86	0.00
32		Benzoic acid	0.175	0.145	17.1	68	0.00
33		4-Chloroaniline	0.412	0.402	2.4	86	0.00
34	C	Hexachlorobutadiene	0.168	0.169	-0.6	89	0.00
35		Caprolactam	0.097	0.091	6.2	82	0.00
36	C	4-Chloro-3-methylphenol	0.288	0.286	0.7	87	0.00
37		2-Methylnaphthalene	0.594	0.579	2.5	86	0.00
38		1-Methylnaphthalene	0.575	0.555	3.5	86	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	87	0.00
40		1,2,4,5-Tetrachlorobenzene	0.616	0.608	1.3	87	0.00
41	P	Hexachlorocyclopentadiene	0.297	0.322	-8.4	91	0.00
42	S	2,4,6-Tribromophenol	0.173	0.192	-11.0	92	0.00
43	C	2,4,6-Trichlorophenol	0.385	0.405	-5.2	90	0.00
44		2,4,5-Trichlorophenol	0.400	0.422	-5.5	89	0.00

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.253	1.226	2.2	90	0.00
46		1,1'-Biphenyl	1.782	1.727	3.1	86	0.00
47		2-Chloronaphthalene	1.313	1.286	2.1	87	0.00
48		2-Nitroaniline	0.337	0.397	-17.8	99	0.00
49		Acenaphthylene	1.907	1.856	2.7	85	0.00
50		Dimethylphthalate	1.416	1.386	2.1	87	0.00
51		2,6-Dinitrotoluene	0.269	0.324	-20.4	99	0.00
52	C	Acenaphthene	1.214	1.207	0.6	88	0.00
53		3-Nitroaniline	0.327	0.374	-14.4	95	0.00
54	P	2,4-Dinitrophenol	0.069	0.107	-55.1#	135	-0.01
55		Dibenzofuran	1.731	1.683	2.8	87	0.00
56	P	4-Nitrophenol	0.250	0.279	-11.6	94	0.00
57		2,4-Dinitrotoluene	0.316	0.421	-33.2#	112	0.00
58		Fluorene	1.260	1.231	2.3	87	0.00
59		2,3,4,6-Tetrachlorophenol	0.317	0.338	-6.6	89	0.00
60		Diethylphthalate	1.399	1.361	2.7	86	0.00
61		4-Chlorophenyl-phenylether	0.656	0.648	1.2	87	0.00
62		4-Nitroaniline	0.310	0.365	-17.7	98	0.00
63		Azobenzene	1.461	1.353	7.4	83	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
65		4,6-Dinitro-2-methylphenol	0.064	0.095	-48.4#	124	-0.01
66	c	n-Nitrosodiphenylamine	0.666	0.653	2.0	86	0.00
67		4-Bromophenyl-phenylether	0.217	0.215	0.9	86	0.00
68		Hexachlorobenzene	0.231	0.229	0.9	88	-0.01
69		Atrazine	0.208	0.213	-2.4	88	0.00
70	C	Pentachlorophenol	0.132	0.146	-10.6	90	0.00
71		Phenanthrene	1.035	1.007	2.7	86	0.00
72		Anthracene	1.041	1.017	2.3	86	-0.01
73		Carbazole	1.023	0.974	4.8	84	0.00
74		Di-n-butylphthalate	1.142	1.113	2.5	85	0.00
75	C	Fluoranthene	1.034	0.996	3.7	86	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	83	0.00
77		Benzidine	0.767	0.783	-2.1	81	0.00
78		Pyrene	1.469	1.529	-4.1	85	-0.01
79	S	Terphenyl-d14	0.952	0.970	-1.9	85	0.00
80		Butylbenzylphthalate	0.594	0.640	-7.7	85	0.00
81		Benzo(a)anthracene	1.178	1.160	1.5	83	0.00
82		3,3'-Dichlorobenzidine	0.413	0.404	2.2	79	0.00
83		Chrysene	1.121	1.100	1.9	82	0.00
84		Bis(2-ethylhexyl)phthalate	0.781	0.804	-2.9	83	-0.01
85	c	Di-n-octyl phthalate	1.102	1.177	-6.8	87	0.00
86		Indeno(1,2,3-cd)pyrene	0.903	0.972	-7.6	95	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	83	-0.01

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88	Benzo(b)fluoranthene	1.153	1.173	-1.7	80	0.00
89	Benzo(k)fluoranthene	1.137	1.141	-0.4	84	-0.01
90 C	Benzo(a)pyrene	1.060	1.079	-1.8	83	-0.01
91	Dibenzo(a,h)anthracene	0.940	1.081	-15.0	93	-0.02
92	Benzo(g,h,i)perylene	0.950	1.111	-16.9	94	-0.02

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