Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022119\
Data File : BM018892.D
Acq On : 21 Feb 2019 12:56
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Feb 22 12:58:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 11 16:02:12 2019
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	124	-0.02
2		1,4-Dioxane	0.532	0.504	5.3	118	0.00
3		Pyridine	1.450	1.574	-8.6	123	0.00
4		n-Nitrosodimethylamine	0.369	0.423	-14.6	136	-0.01
5	S	2-Fluorophenol	1.221	1.263	-3.4	126	-0.02
6		Aniline	2.107	2.267	-7.6	130	-0.01
7	S	Phenol-d6	1.720	1.882	-9.4	132	-0.01
8		2-Chlorophenol	1.340	1.444	-7.8	131	-0.02
9		Benzaldehyde	0.941	1.044	-10.9	138	-0.02
10	C	Phenol	1.809	1.961	-8.4	132	-0.01
11		bis(2-Chloroethyl)ether	1.380	1.488	-7.8	131	-0.01
12		1,3-Dichlorobenzene	1.507	1.570	-4.2	128	-0.01
13	C	1,4-Dichlorobenzene	1.534	1.608	-4.8	130	-0.02
14		1,2-Dichlorobenzene	1.474	1.554	-5.4	130	-0.02
15		Benzyl Alcohol	1.366	1.529	-11.9	132	-0.01
16		2,2'-oxybis(1-Chloropropane	1.336	1.457	-9.1	137	-0.02
17		2-Methylphenol	1.152	1.254	-8.9	132	-0.02
18		Hexachloroethane	0.560	0.590	-5.4	129	-0.02
19	P	n-Nitroso-di-n-propylamine	1.331	1.473	-10.7	132	-0.02
20		3+4-Methylphenols	1.590	1.761	-10.8	132	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	127	-0.02
22		Acetophenone	0.549	0.576	-4.9	133	-0.02
23	S	Nitrobenzene-d5	0.433	0.463	-6.9	134	-0.02
24		Nitrobenzene	0.449	0.473	-5.3	132	-0.02
25		Isophorone	0.690	0.736	-6.7	132	-0.02
26	C	2-Nitrophenol	0.179	0.198	-10.6	135	-0.02
27		2,4-Dimethylphenol	0.261	0.277	-6.1	133	-0.01
28		bis(2-Chloroethoxy)methane	0.446	0.468	-4.9	131	-0.02
29	C	2,4-Dichlorophenol	0.300	0.327	-9.0	134	-0.02
30		1,2,4-Trichlorobenzene	0.346	0.359	-3.8	132	-0.02
31		Naphthalene	0.975	1.013	-3.9	133	-0.02
32		Benzoic acid	0.228	0.204	10.5	113	0.00
33		4-Chloroaniline	0.436	0.465	-6.7	133	-0.01
34	C	Hexachlorobutadiene	0.201	0.202	-0.5	130	-0.02
35		Caprolactam	0.122	0.129	-5.7	130	0.00
36	C	4-Chloro-3-methylphenol	0.362	0.391	-8.0	134	-0.01
37		2-Methylnaphthalene	0.687	0.721	-4.9	133	-0.02
38		1-Methylnaphthalene	0.672	0.700	-4.2	133	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	126	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.640	0.661	-3.3	131	-0.02
41	P	Hexachlorocyclopentadiene	0.327	0.286	12.5	109	-0.02
42	S	2,4,6-Tribromophenol	0.288	0.315	-9.4	134	-0.01
43	C	2,4,6-Trichlorophenol	0.424	0.459	-8.3	132	-0.01
44		2,4,5-Trichlorophenol	0.444	0.476	-7.2	131	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.507	1.571	-4.2	133	-0.02
46		1,1'-Biphenyl	1.722	1.805	-4.8	134	-0.01
47		2-Chloronaphthalene	1.330	1.410	-6.0	134	-0.02
48		2-Nitroaniline	0.442	0.502	-13.6	136	-0.01
49		Acenaphthylene	1.980	2.092	-5.7	132	-0.01
50		Dimethylphthalate	1.667	1.718	-3.1	130	-0.01
51		2,6-Dinitrotoluene	0.361	0.387	-7.2	130	-0.01
52	C	Acenaphthene	1.214	1.272	-4.8	132	-0.01
53		3-Nitroaniline	0.408	0.422	-3.4	129	-0.01
54	P	2,4-Dinitrophenol	0.200	0.164	18.0	103	-0.01
55		Dibenzofuran	1.995	2.090	-4.8	133	-0.02
56	P	4-Nitrophenol	0.326	0.342	-4.9	130	-0.01
57		2,4-Dinitrotoluene	0.497	0.547	-10.1	132	-0.01
58		Fluorene	1.527	1.614	-5.7	134	-0.01
59		2,3,4,6-Tetrachlorophenol	0.392	0.408	-4.1	128	-0.01
60		Diethylphthalate	1.719	1.775	-3.3	131	-0.01
61		4-Chlorophenyl-phenylether	0.856	0.888	-3.7	131	-0.01
62		4-Nitroaniline	0.400	0.394	1.5	122	-0.01
63		Azobenzene	1.844	1.995	-8.2	134	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	124	-0.01
65		4,6-Dinitro-2-methylphenol	0.140	0.126	10.0	111	0.00
66	c	n-Nitrosodiphenylamine	0.632	0.679	-7.4	132	-0.01
67		4-Bromophenyl-phenylether	0.224	0.237	-5.8	131	-0.02
68		Hexachlorobenzene	0.260	0.275	-5.8	132	-0.01
69		Atrazine	0.204	0.190	6.9	115	-0.01
70	C	Pentachlorophenol	0.168	0.178	-6.0	124	-0.01
71		Phenanthrene	1.075	1.133	-5.4	133	-0.01
72		Anthracene	1.046	1.124	-7.5	132	-0.02
73		Carbazole	1.093	1.143	-4.6	128	-0.01
74		Di-n-butylphthalate	1.257	1.378	-9.6	132	-0.01
75	C	Fluoranthene	1.242	1.305	-5.1	129	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	112	-0.01
77		Benzidine	0.450	0.382	15.1	77	0.00
78		Pyrene	1.162	1.339	-15.2	128	-0.01
79	S	Terphenyl-d14	0.970	1.137	-17.2	127	0.00
80		Butylbenzylphthalate	0.529	0.650	-22.9	130	-0.01
81		Benzo(a)anthracene	1.166	1.245	-6.8	118	0.00
82		3,3'-Dichlorobenzidine	0.461	0.473	-2.6	111	-0.01
83		Chrysene	1.117	1.183	-5.9	117	0.00
84		Bis(2-ethylhexyl)phthalate	0.775	0.949	-22.5	132	-0.01
85	c	Di-n-octyl phthalate	1.302	1.580	-21.4#	129	-0.01
86		Indeno(1,2,3-cd)pyrene	1.290	1.072	16.9	85	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	91	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.144	1.262	-10.3	101	-0.01
89	Benzo(k)fluoranthene	1.139	1.295	-13.7	105	-0.02
90 C	Benzo(a)pyrene	1.084	1.161	-7.1	97	-0.02
91	Dibenzo(a,h)anthracene	1.074	1.071	0.3	86	-0.02
92	Benzo(g,h,i)perylene	1.000	0.977	2.3	83	-0.03

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