Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082420\
Data File : BM027194.D
Acq On : 24 Aug 2020 14:09
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 24 15:38:00 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 24 15:18:39 2020
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	158#	0.00
2		1,4-Dioxane	0.490	0.428	12.7	151#	0.00
3		Pyridine	1.376	1.278	7.1	153#	0.00
4		n-Nitrosodimethylamine	0.600	0.485	19.2	137	0.00
5	S	2-Fluorophenol	1.026	1.043	-1.7	165#	0.00
6		Aniline	1.600	1.556	2.8	217#	0.00
7	S	Phenol-d6	1.409	1.405	0.3	218#	0.00
8		2-Chlorophenol	1.068	1.038	2.8	213#	0.00
9		Benzaldehyde	0.964	0.945	2.0	220#	0.00
10	C	Phenol	1.342	1.298	3.3	216#	0.00
11		bis(2-Chloroethyl)ether	1.097	1.087	0.9	222#	0.00
12		1,3-Dichlorobenzene	1.494	1.376	7.9	161#	0.00
13	C	1,4-Dichlorobenzene	1.508	1.385	8.2	159#	0.00
14		1,2-Dichlorobenzene	1.430	1.308	8.5	158#	0.00
15		Benzyl Alcohol	1.280	1.166	8.9	153#	0.00
16		2,2'-oxybis(1-Chloropropane	0.880	0.859	2.4	164#	0.00
17		2-Methylphenol	0.929	0.884	4.8	161#	0.00
18		Hexachloroethane	0.612	0.544	11.1	155#	0.00
19	P	n-Nitroso-di-n-propylamine	1.081	0.992	8.2	153#	0.00
20		3+4-Methylphenols	1.243	1.195	3.9	160#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	155#	0.00
22		Acetophenone	0.538	0.490	8.9	155#	0.00
23	S	Nitrobenzene-d5	0.485	0.430	11.3	149	0.00
24		Nitrobenzene	0.491	0.422	14.1	146	0.00
25		Isophorone	0.710	0.655	7.7	152#	0.00
26	C	2-Nitrophenol	0.153	0.164	-7.2	177#	0.00
27		2,4-Dimethylphenol	0.238	0.225	5.5	157#	0.00
28		bis(2-Chloroethoxy)methane	0.433	0.415	4.2	158#	0.00
29	C	2,4-Dichlorophenol	0.332	0.315	5.1	157#	0.00
30		1,2,4-Trichlorobenzene	0.435	0.396	9.0	156#	0.00
31		Naphthalene	1.014	0.915	9.8	153#	0.00
32		Benzoic acid	0.170	0.191	-12.4	167#	0.00
33		4-Chloroaniline	0.419	0.396	5.5	154#	0.00
34	C	Hexachlorobutadiene	0.314	0.275	12.4	156#	0.00
35		Caprolactam	0.090	0.095	-5.6	163#	0.00
36	C	4-Chloro-3-methylphenol	0.343	0.322	6.1	150	0.00
37		2-Methylnaphthalene	0.768	0.694	9.6	150#	0.00
38		1-Methylnaphthalene	0.739	0.662	10.4	150	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	149	0.00
40		1,2,4,5-Tetrachlorobenzene	0.704	0.637	9.5	154#	0.00
41	P	Hexachlorocyclopentadiene	0.389	0.349	10.3	147	0.00
42	S	2,4,6-Tribromophenol	0.324	0.305	5.9	149	0.00
43	C	2,4,6-Trichlorophenol	0.435	0.416	4.4	156#	0.00
44		2,4,5-Trichlorophenol	0.461	0.430	6.7	154#	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.523	1.360	10.7	150	0.00
46		1,1'-Biphenyl	1.499	1.343	10.4	148	0.00
47		2-Chloronaphthalene	1.278	1.147	10.3	150	0.00
48		2-Nitroaniline	0.408	0.380	6.9	143	0.00
49		Acenaphthylene	1.777	1.657	6.8	148	0.00
50		Dimethylphthalate	1.624	1.465	9.8	145	0.00
51		2,6-Dinitrotoluene	0.323	0.310	4.0	148	0.00
52	C	Acenaphthene	1.128	1.042	7.6	147	0.00
53		3-Nitroaniline	0.311	0.312	-0.3	145	0.00
54	P	2,4-Dinitrophenol	0.169	0.168	0.6	154#	0.00
55		Dibenzofuran	1.880	1.690	10.1	143	0.00
56	P	4-Nitrophenol	0.248	0.256	-3.2	155#	0.00
57		2,4-Dinitrotoluene	0.449	0.437	2.7	146	0.00
58		Fluorene	1.552	1.416	8.8	144	0.00
59		2,3,4,6-Tetrachlorophenol	0.452	0.427	5.5	150#	0.00
60		Diethylphthalate	1.611	1.498	7.0	145	0.00
61		4-Chlorophenyl-phenylether	0.891	0.795	10.8	146	0.00
62		4-Nitroaniline	0.349	0.352	-0.9	147	0.00
63		Azobenzene	1.653	1.472	10.9	138	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	144	0.00
65		4,6-Dinitro-2-methylphenol	0.098	0.101	-3.1	158#	0.00
66	c	n-Nitrosodiphenylamine	0.568	0.528	7.0	144	0.00
67		4-Bromophenyl-phenylether	0.241	0.223	7.5	146	0.00
68		Hexachlorobenzene	0.287	0.260	9.4	145	0.00
69		Atrazine	0.215	0.205	4.7	147	0.00
70	C	Pentachlorophenol	0.155	0.157	-1.3	155#	0.00
71		Phenanthrene	1.112	1.010	9.2	140	0.00
72		Anthracene	1.063	0.985	7.3	142	0.00
73		Carbazole	0.978	0.918	6.1	143	0.00
74		Di-n-butylphthalate	1.129	1.142	-1.2	150#	0.00
75	C	Fluoranthene	1.389	1.126	18.9	125	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	129	0.00
77		Benzidine	0.347	0.408	-17.6	138	0.00
78		Pyrene	1.217	1.137	6.6	125	0.00
79	S	Terphenyl-d14	1.071	1.002	6.4	127	0.00
80		Butylbenzylphthalate	0.445	0.442	0.7	126	0.00
81		Benzo(a)anthracene	1.296	1.204	7.1	126	0.00
82		3,3'-Dichlorobenzidine	0.452	0.467	-3.3	136	0.00
83		Chrysene	1.259	1.158	8.0	126	0.00
84		Bis(2-ethylhexyl)phthalate	0.633	0.638	-0.8	126	0.00
85	c	Di-n-octyl phthalate	1.044	1.071	-2.6	129	0.00
86	I	Perylene-d12	1.000	1.000	0.0	135	0.00
87		Indeno(1,2,3-cd)pyrene	1.565	1.271	18.8	122	0.00

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88	Benzo(b)fluoranthene	1.334	1.161	13.0	121	0.00
89	Benzo(k)fluoranthene	1.306	1.147	12.2	126	0.00
90 C	Benzo(a)pyrene	1.218	1.077	11.6	125	0.00
91	Dibenzo(a,h)anthracene	1.303	1.055	19.0	123	0.00
92	Benzo(g,h,i)perylene	1.255	1.010	19.5	123	0.00

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