Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM013020\
Data File : BM024659.D
Acq On : 30 Jan 2020 15:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jan 30 23:36:21 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 23 14:57:07 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	77	-0.01
2		1,4-Dioxane	0.416	0.417	-0.2	76	-0.01
3		Pyridine	1.206	1.215	-0.7	78	-0.01
4		n-Nitrosodimethylamine	0.535	0.430	19.6	60	-0.01
5	S	2-Fluorophenol	1.052	1.153	-9.6	82	-0.01
6		Aniline	1.764	1.774	-0.6	75	-0.02
7	S	Phenol-d6	1.449	1.481	-2.2	76	-0.01
8		2-Chlorophenol	1.197	1.315	-9.9	82	-0.01
9		Benzaldehyde	0.856	0.791	7.6	72	-0.02
10	C	Phenol	1.444	1.458	-1.0	75	-0.01
11		bis(2-Chloroethyl)ether	1.156	1.121	3.0	71	-0.01
12		1,3-Dichlorobenzene	1.471	1.660	-12.8	84	-0.01
13	C	1,4-Dichlorobenzene	1.492	1.671	-12.0	83	-0.02
14		1,2-Dichlorobenzene	1.440	1.614	-12.1	84	-0.02
15		Benzyl Alcohol	1.194	1.119	6.3	70	-0.01
16		2,2'-oxybis(1-Chloropropane	1.058	0.897	15.2	62	-0.02
17		2-Methylphenol	1.018	1.030	-1.2	75	0.00
18		Hexachloroethane	0.576	0.580	-0.7	75	-0.02
19	P	n-Nitroso-di-n-propylamine	1.088	0.930	14.5	64	-0.02
20		3+4-Methylphenols	1.409	1.398	0.8	73	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	69	-0.01
22		Acetophenone	0.516	0.531	-2.9	70	-0.01
23	S	Nitrobenzene-d5	0.408	0.392	3.9	65	-0.02
24		Nitrobenzene	0.391	0.377	3.6	65	-0.02
25		Isophorone	0.693	0.657	5.2	64	-0.02
26	C	2-Nitrophenol	0.181	0.208	-14.9	77	-0.01
27		2,4-Dimethylphenol	0.246	0.266	-8.1	74	-0.01
28		bis(2-Chloroethoxy)methane	0.423	0.426	-0.7	68	-0.01
29	C	2,4-Dichlorophenol	0.332	0.382	-15.1	77	-0.01
30		1,2,4-Trichlorobenzene	0.401	0.466	-16.2	80	-0.01
31		Naphthalene	1.010	1.103	-9.2	74	-0.01
32		Benzoic acid	0.225	0.249	-10.7	73	-0.03
33		4-Chloroaniline	0.443	0.476	-7.4	72	-0.01
34	C	Hexachlorobutadiene	0.277	0.316	-14.1	77	-0.01
35		Caprolactam	0.107	0.113	-5.6	71	-0.02
36	C	4-Chloro-3-methylphenol	0.353	0.351	0.6	67	-0.01
37		2-Methylnaphthalene	0.780	0.831	-6.5	72	-0.02
38		1-Methylnaphthalene	0.742	0.786	-5.9	72	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	68	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.714	0.807	-13.0	75	-0.01
41	P	Hexachlorocyclopentadiene	0.400	0.365	8.8	59	-0.02
42	S	2,4,6-Tribromophenol	0.325	0.376	-15.7	77	-0.01
43	C	2,4,6-Trichlorophenol	0.454	0.518	-14.1	74	-0.01
44		2,4,5-Trichlorophenol	0.464	0.525	-13.1	74	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.471	1.615	-9.8	72	-0.01
46		1,1'-Biphenyl	1.514	1.666	-10.0	72	-0.02
47		2-Chloronaphthalene	1.222	1.372	-12.3	74	-0.01
48		2-Nitroaniline	0.373	0.332	11.0	58	-0.01
49		Acenaphthylene	1.829	2.008	-9.8	72	-0.01
50		Dimethylphthalate	1.617	1.716	-6.1	70	-0.01
51		2,6-Dinitrotoluene	0.344	0.373	-8.4	71	-0.01
52	C	Acenaphthene	1.135	1.234	-8.7	72	-0.01
53		3-Nitroaniline	0.357	0.386	-8.1	70	-0.01
54	P	2,4-Dinitrophenol	0.206	0.220	-6.8	70	0.00
55		Dibenzofuran	1.851	2.035	-9.9	73	-0.01
56	P	4-Nitrophenol	0.279	0.298	-6.8	69	0.00
57		2,4-Dinitrotoluene	0.493	0.521	-5.7	69	0.00
58		Fluorene	1.556	1.640	-5.4	69	-0.01
59		2,3,4,6-Tetrachlorophenol	0.470	0.515	-9.6	72	-0.01
60		Diethylphthalate	1.642	1.666	-1.5	66	-0.01
61		4-Chlorophenyl-phenylether	0.897	0.941	-4.9	69	-0.01
62		4-Nitroaniline	0.394	0.418	-6.1	68	-0.01
63		Azobenzene	1.401	1.235	11.8	58	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	66	-0.01
65		4,6-Dinitro-2-methylphenol	0.118	0.132	-11.9	71	-0.01
66	c	n-Nitrosodiphenylamine	0.574	0.630	-9.8	71	-0.01
67		4-Bromophenyl-phenylether	0.249	0.274	-10.0	71	-0.01
68		Hexachlorobenzene	0.285	0.330	-15.8	75	-0.01
69		Atrazine	0.221	0.213	3.6	61	-0.01
70	C	Pentachlorophenol	0.174	0.187	-7.5	69	0.00
71		Phenanthrene	1.083	1.198	-10.6	71	-0.01
72		Anthracene	1.060	1.176	-10.9	72	-0.01
73		Carbazole	0.965	1.083	-12.2	72	-0.01
74		Di-n-butylphthalate	1.233	1.273	-3.2	66	-0.01
75	C	Fluoranthene	1.352	1.523	-12.6	73	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	70	-0.01
77		Benzidine	0.466	0.592	-27.0#	86	0.00
78		Pyrene	1.318	1.430	-8.5	73	0.00
79	S	Terphenyl-d14	1.071	1.190	-11.1	70	-0.01
80		Butylbenzylphthalate	0.553	0.563	-1.8	68	0.00
81		Benzo(a)anthracene	1.388	1.514	-9.1	74	0.00
82		3,3'-Dichlorobenzidine	0.499	0.612	-22.6	80	0.00
83		Chrysene	1.325	1.466	-10.6	75	0.00
84		Bis(2-ethylhexyl)phthalate	0.816	0.802	1.7	66	0.00
85	c	Di-n-octyl phthalate	1.349	1.393	-3.3	68	0.00
86		Indeno(1,2,3-cd)pyrene	1.444	1.863	-29.0#	84	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	74	0.00

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88	Benzo(b)fluoranthene	1.297	1.360	-4.9	75	0.00
89	Benzo(k)fluoranthene	1.266	1.337	-5.6	77	-0.01
90 C	Benzo(a)pyrene	1.179	1.285	-9.0	79	-0.01
91	Dibenzo(a,h)anthracene	1.119	1.330	-18.9	84	-0.01
92	Benzo(g,h,i)perylene	1.016	1.281	-26.1#	88	-0.02

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