Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
Data File : BM020127.D
Acq On : 04 May 2019 00:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 04 06:54:07 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 01 03:58:37 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	209#	0.00
2		1,4-Dioxane	0.600	0.558	7.0	209#	0.00
3		Pyridine	1.535	1.561	-1.7	207#	0.00
4		n-Nitrosodimethylamine	0.492	0.521	-5.9	232#	0.00
5	S	2-Fluorophenol	1.224	1.264	-3.3	225#	0.00
6		Aniline	2.104	2.273	-8.0	236#	0.00
7	S	Phenol-d6	1.672	1.827	-9.3	237#	0.00
8		2-Chlorophenol	1.332	1.423	-6.8	229#	0.00
9		Benzaldehyde	1.265	1.307	-3.3	218#	0.00
10	C	Phenol	1.799	1.959	-8.9	233#	0.00
11		bis(2-Chloroethyl)ether	1.308	1.386	-6.0	224#	0.00
12		1,3-Dichlorobenzene	1.550	1.620	-4.5	226#	0.00
13	C	1,4-Dichlorobenzene	1.566	1.632	-4.2	224#	0.00
14		1,2-Dichlorobenzene	1.492	1.567	-5.0	229#	0.00
15		Benzyl Alcohol	1.612	1.731	-7.4	228#	0.00
16		2,2'-oxybis(1-Chloropropane	1.084	1.103	-1.8	210#	0.00
17		2-Methylphenol	1.158	1.245	-7.5	228#	0.00
18		Hexachloroethane	0.653	0.664	-1.7	222#	0.00
19	P	n-Nitroso-di-n-propylamine	1.430	1.518	-6.2	222#	0.00
20		3+4-Methylphenols	1.540	1.688	-9.6	232#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	208#	0.00
22		Acetophenone	0.547	0.570	-4.2	225#	0.00
23	S	Nitrobenzene-d5	0.507	0.530	-4.5	225#	0.00
24		Nitrobenzene	0.525	0.540	-2.9	222#	0.00
25		Isophorone	0.874	0.933	-6.8	223#	0.00
26	C	2-Nitrophenol	0.159	0.198	-24.5#	264#	0.00
27		2,4-Dimethylphenol	0.264	0.285	-8.0	233#	0.00
28		bis(2-Chloroethoxy)methane	0.476	0.512	-7.6	226#	0.00
29	C	2,4-Dichlorophenol	0.333	0.371	-11.4	239#	0.00
30		1,2,4-Trichlorobenzene	0.410	0.437	-6.6	230#	0.00
31		Naphthalene	1.038	1.100	-6.0	228#	0.00
32		Benzoic acid	0.175	0.213	-21.7	255#	0.04
33		4-Chloroaniline	0.427	0.461	-8.0	227#	0.00
34	C	Hexachlorobutadiene	0.290	0.306	-5.5	226#	-0.01
35		Caprolactam	0.100	0.116	-16.0	237#	0.02
36	C	4-Chloro-3-methylphenol	0.391	0.426	-9.0	226#	0.00
37		2-Methylnaphthalene	0.757	0.826	-9.1	231#	0.00
38		1-Methylnaphthalene	0.740	0.799	-8.0	229#	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	202#	0.00
40		1,2,4,5-Tetrachlorobenzene	0.782	0.836	-6.9	228#	0.00
41	P	Hexachlorocyclopentadiene	0.461	0.273	40.8#	127	0.00
42	S	2,4,6-Tribromophenol	0.310	0.355	-14.5	233#	0.00
43	C	2,4,6-Trichlorophenol	0.445	0.518	-16.4	242#	0.00
44		2,4,5-Trichlorophenol	0.497	0.570	-14.7	242#	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.626	1.739	-6.9	228#	0.00
46		1,1'-Biphenyl	1.646	1.732	-5.2	222#	0.00
47		2-Chloronaphthalene	1.314	1.394	-6.1	228#	0.00
48		2-Nitroaniline	0.468	0.511	-9.2	224#	0.00
49		Acenaphthylene	2.103	2.242	-6.6	223#	0.00
50		Dimethylphthalate	1.700	1.762	-3.6	214#	0.00
51		2,6-Dinitrotoluene	0.358	0.390	-8.9	225#	0.00
52	C	Acenaphthene	1.206	1.279	-6.1	221#	0.00
53		3-Nitroaniline	0.332	0.365	-9.9	226#	0.00
54	P	2,4-Dinitrophenol	0.151	0.149	1.3	211#	0.00
55		Dibenzofuran	2.035	2.132	-4.8	220#	0.00
56	P	4-Nitrophenol	0.284	0.315	-10.9	224#	0.00
57		2,4-Dinitrotoluene	0.486	0.544	-11.9	227#	0.00
58		Fluorene	1.671	1.781	-6.6	223#	0.00
59		2,3,4,6-Tetrachlorophenol	0.472	0.537	-13.8	237#	0.00
60		Diethylphthalate	1.713	1.784	-4.1	211#	0.00
61		4-Chlorophenyl-phenylether	0.947	1.016	-7.3	224#	0.00
62		4-Nitroaniline	0.367	0.400	-9.0	220#	0.00
63		Azobenzene	2.020	2.037	-0.8	208#	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	199#	0.00
65		4,6-Dinitro-2-methylphenol	0.094	0.092	2.1	209#	0.00
66	c	n-Nitrosodiphenylamine	0.589	0.630	-7.0	219#	0.00
67		4-Bromophenyl-phenylether	0.245	0.267	-9.0	223#	0.00
68		Hexachlorobenzene	0.282	0.302	-7.1	219#	0.00
69		Atrazine	0.230	0.243	-5.7	213#	0.00
70	C	Pentachlorophenol	0.161	0.194	-20.5#	243#	0.00
71		Phenanthrene	1.104	1.172	-6.2	218#	0.00
72		Anthracene	1.104	1.170	-6.0	218#	0.00
73		Carbazole	0.996	1.045	-4.9	215#	0.00
74		Di-n-butylphthalate	1.199	1.284	-7.1	214#	0.00
75	C	Fluoranthene	1.397	1.452	-3.9	213#	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	185#	0.00
77		Benzidine	0.639	0.682	-6.7	201#	0.00
78		Pyrene	1.343	1.485	-10.6	212#	0.00
79	S	Terphenyl-d14	1.147	1.263	-10.1	210#	0.00
80		Butylbenzylphthalate	0.488	0.578	-18.4	219#	-0.01
81		Benzo(a)anthracene	1.403	1.523	-8.6	209#	0.00
82		3,3'-Dichlorobenzidine	0.535	0.607	-13.5	213#	0.00
83		Chrysene	1.360	1.446	-6.3	202#	0.00
84		Bis(2-ethylhexyl)phthalate	0.758	0.871	-14.9	210#	-0.01
85	c	Di-n-octyl phthalate	1.259	1.455	-15.6	214#	-0.02
86		Indeno(1,2,3-cd)pyrene	1.618	1.748	-8.0	210#	0.00
87	I	Perylene-d12	1.000	1.000	0.0	191#	-0.01

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88	Benzo(b)fluoranthene	1.321	1.374	-4.0	203#	0.00
89	Benzo(k)fluoranthene	1.205	1.261	-4.6	205#	0.00
90 C	Benzo(a)pyrene	1.200	1.271	-5.9	207#	0.00
91	Dibenzo(a,h)anthracene	1.248	1.317	-5.5	210#	0.00
92	Benzo(g,h,i)perylene	1.184	1.255	-6.0	210#	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 2