Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040121\
Data File : BM029308.D
Acq On : 01 Apr 2021 16:24
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040121

Quant Time: Apr 01 17:17:31 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 01 16:52:48 2021
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	104	0.00
2		1,4-Dioxane	0.526	0.502	4.6	99	0.00
3		Pyridine	1.331	1.344	-1.0	98	0.00
4		n-Nitrosodimethylamine	0.635	0.662	-4.3	101	0.00
5	S	2-Fluorophenol	1.211	1.192	1.6	102	0.00
6		Aniline	1.867	1.865	0.1	101	0.00
7	S	Phenol-d6	1.710	1.706	0.2	102	0.00
8		2-Chlorophenol	1.341	1.340	0.1	103	0.00
9		Benzaldehyde	1.076	1.044	3.0	104	0.00
10	C	Phenol	1.682	1.676	0.4	101	0.00
11		bis(2-Chloroethyl)ether	1.273	1.275	-0.2	101	0.00
12		1,3-Dichlorobenzene	1.459	1.430	2.0	100	0.00
13	C	1,4-Dichlorobenzene	1.484	1.447	2.5	101	0.00
14		1,2-Dichlorobenzene	1.428	1.407	1.5	101	0.00
15		Benzyl Alcohol	1.261	1.293	-2.5	102	0.00
16		2,2'-oxybis(1-Chloropropane	1.909	1.927	-0.9	102	0.00
17		2-Methylphenol	1.096	1.106	-0.9	101	0.00
18		Hexachloroethane	0.581	0.565	2.8	101	0.00
19	P	n-Nitroso-di-n-propylamine	1.156	1.187	-2.7	101	0.00
20		3+4-Methylphenols	1.485	1.502	-1.1	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
22		Acetophenone	0.517	0.531	-2.7	101	0.00
23	S	Nitrobenzene-d5	0.436	0.446	-2.3	101	0.00
24		Nitrobenzene	0.448	0.458	-2.2	102	0.00
25		Isophorone	0.781	0.812	-4.0	101	0.00
26	C	2-Nitrophenol	0.179	0.186	-3.9	102	0.00
27		2,4-Dimethylphenol	0.281	0.289	-2.8	101	0.00
28		bis(2-Chloroethoxy)methane	0.403	0.411	-2.0	100	0.00
29	C	2,4-Dichlorophenol	0.311	0.318	-2.3	101	0.00
30		1,2,4-Trichlorobenzene	0.336	0.338	-0.6	102	0.00
31		Naphthalene	1.054	1.063	-0.9	101	0.00
32		Benzoic acid	0.212	0.225	-6.1	104	0.00
33		4-Chloroaniline	0.425	0.440	-3.5	100	0.00
34	C	Hexachlorobutadiene	0.198	0.200	-1.0	103	0.00
35		Caprolactam	0.098	0.107	-9.2	102	0.00
36	C	4-Chloro-3-methylphenol	0.335	0.347	-3.6	100	0.00
37		2-Methylnaphthalene	0.743	0.765	-3.0	101	0.00
38		1-Methylnaphthalene	0.703	0.718	-2.1	101	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
40		1,2,4,5-Tetrachlorobenzene	0.561	0.565	-0.7	102	0.00
41	P	Hexachlorocyclopentadiene	0.325	0.329	-1.2	100	0.00
42	S	2,4,6-Tribromophenol	0.200	0.215	-7.5	104	0.00
43	C	2,4,6-Trichlorophenol	0.398	0.412	-3.5	101	0.00
44		2,4,5-Trichlorophenol	0.429	0.446	-4.0	102	0.00
45	S	2-Fluorobiphenyl	1.436	1.452	-1.1	102	0.00
46		1,1'-Biphenyl	1.548	1.564	-1.0	101	0.00
47		2-Chloronaphthalene	1.198	1.208	-0.8	101	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.400	0.424	-6.0	102	0.00
49		Acenaphthylene	1.868	1.929	-3.3	102	0.00
50		Dimethylphthalate	1.444	1.476	-2.2	101	0.00
51		2,6-Dinitrotoluene	0.321	0.336	-4.7	101	0.00
52	C	Acenaphthene	1.155	1.177	-1.9	101	0.00
53		3-Nitroaniline	0.328	0.352	-7.3	102	0.00
54	P	2,4-Dinitrophenol	0.169	0.189	-11.8	102	0.00
55		Dibenzofuran	1.811	1.854	-2.4	102	0.00
56	P	4-Nitrophenol	0.280	0.310	-10.7	101	0.00
57		2,4-Dinitrotoluene	0.426	0.460	-8.0	102	0.00
58		Fluorene	1.486	1.523	-2.5	101	0.00
59		2,3,4,6-Tetrachlorophenol	0.330	0.348	-5.5	102	0.00
60		Diethylphthalate	1.475	1.523	-3.3	101	0.00
61		4-Chlorophenyl-phenylether	0.680	0.702	-3.2	102	0.00
62		4-Nitroaniline	0.345	0.375	-8.7	102	0.00
63		Azobenzene	1.697	1.768	-4.2	101	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
65		4,6-Dinitro-2-methylphenol	0.128	0.132	-3.1	101	0.00
66	c	n-Nitrosodiphenylamine	0.661	0.661	0.0	100	0.00
67		4-Bromophenyl-phenylether	0.199	0.201	-1.0	101	0.00
68		Hexachlorobenzene	0.218	0.218	0.0	103	0.00
69		Atrazine	0.220	0.227	-3.2	102	0.00
70	C	Pentachlorophenol	0.102	0.122	-19.6	112	0.00
71		Phenanthrene	1.157	1.164	-0.6	102	0.00
72		Anthracene	1.142	1.169	-2.4	103	0.00
73		Carbazole	1.115	1.145	-2.7	103	0.00
74		Di-n-butylphthalate	1.279	1.315	-2.8	100	0.00
75	C	Fluoranthene	1.289	1.314	-1.9	102	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
77		Benzidine	0.649	0.708	-9.1	104	0.00
78		Pyrene	1.390	1.380	0.7	102	0.00
79	S	Terphenyl-d14	1.053	1.060	-0.7	103	0.00
80		Butylbenzylphthalate	0.626	0.629	-0.5	101	0.00
81		Benzo(a)anthracene	1.335	1.339	-0.3	102	0.00
82		3,3'-Dichlorobenzidine	0.448	0.451	-0.7	101	0.00
83		Chrysene	1.307	1.308	-0.1	102	0.00
84		Bis(2-ethylhexyl)phthalate	0.963	0.977	-1.5	102	0.00
85	c	Di-n-octyl phthalate	1.652	1.658	-0.4	100	0.00
86	I	Perylene-d12	1.000	1.000	0.0	101	0.00
87		Indeno(1,2,3-cd)pyrene	1.474	1.517	-2.9	103	0.00
88		Benzo(b)fluoranthene	1.313	1.361	-3.7	106	0.00
89		Benzo(k)fluoranthene	1.256	1.263	-0.6	98	0.00
90	C	Benzo(a)pyrene	1.201	1.223	-1.8	102	0.00
91		Dibenzo(a,h)anthracene	1.266	1.299	-2.6	102	0.00
92		Benzo(g,h,i)perylene	1.220	1.231	-0.9	102	0.00

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