Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091823\
Data File : BM041897.D
Acq On : 18 Sep 2023 16:12
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM091823

Quant Time: Sep 19 04:51:56 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM091823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 19 04:49:34 2023
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	103	0.00
2		1,4-Dioxane	0.490	0.489	0.2	106	0.00
3		Pyridine	1.521	1.515	0.4	103	0.00
4		n-Nitrosodimethylamine	0.723	0.737	-1.9	105	0.00
5	S	2-Fluorophenol	1.201	1.204	-0.2	105	0.00
6		Aniline	2.209	2.217	-0.4	104	0.00
7	S	Phenol-d6	1.667	1.689	-1.3	104	0.00
8		2-Chlorophenol	1.337	1.342	-0.4	104	0.00
9		Benzaldehyde	0.943	0.943	0.0	107	0.00
10	C	Phenol	1.711	1.730	-1.1	105	0.00
11		bis(2-Chloroethyl)ether	1.428	1.431	-0.2	104	0.00
12		1,3-Dichlorobenzene	1.427	1.417	0.7	103	0.00
13	C	1,4-Dichlorobenzene	1.450	1.439	0.8	103	0.00
14		1,2-Dichlorobenzene	1.394	1.387	0.5	103	0.00
15		Benzyl Alcohol	1.316	1.343	-2.1	104	0.00
16		2,2'-oxybis(1-Chloropropane	2.183	2.185	-0.1	104	0.01
17		2-Methylphenol	1.244	1.263	-1.5	104	0.00
18		Hexachloroethane	0.545	0.538	1.3	102	0.00
19	P	n-Nitroso-di-n-propylamine	1.239	1.272	-2.7	103	0.00
20		3+4-Methylphenols	1.708	1.738	-1.8	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
22		Acetophenone	0.500	0.503	-0.6	104	0.00
23	S	Nitrobenzene-d5	0.393	0.396	-0.8	104	0.00
24		Nitrobenzene	0.390	0.393	-0.8	103	0.00
25		Isophorone	0.775	0.776	-0.1	103	0.00
26	C	2-Nitrophenol	0.176	0.181	-2.8	105	0.00
27		2,4-Dimethylphenol	0.316	0.315	0.3	104	0.00
28		bis(2-Chloroethoxy)methane	0.453	0.454	-0.2	103	0.00
29	C	2,4-Dichlorophenol	0.306	0.311	-1.6	104	0.00
30		1,2,4-Trichlorobenzene	0.314	0.312	0.6	104	0.00
31		Naphthalene	1.007	1.003	0.4	104	0.00
32		Benzoic acid	0.249	0.264	-6.0	105	0.00
33		4-Chloroaniline	0.451	0.452	-0.2	104	0.00
34	C	Hexachlorobutadiene	0.192	0.189	1.6	103	0.00
35		Caprolactam	0.109	0.112	-2.8	104	0.00
36	C	4-Chloro-3-methylphenol	0.341	0.344	-0.9	104	0.00
37		2-Methylnaphthalene	0.740	0.743	-0.4	104	0.00
38		1-Methylnaphthalene	0.695	0.692	0.4	103	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
40		1,2,4,5-Tetrachlorobenzene	0.580	0.586	-1.0	104	0.00
41	P	Hexachlorocyclopentadiene	0.326	0.335	-2.8	102	0.00
42	S	2,4,6-Tribromophenol	0.250	0.251	-0.4	104	0.00
43	C	2,4,6-Trichlorophenol	0.395	0.401	-1.5	104	0.00
44		2,4,5-Trichlorophenol	0.462	0.470	-1.7	104	0.00
45	S	2-Fluorobiphenyl	1.363	1.360	0.2	103	0.00
46		1,1'-Biphenyl	1.476	1.480	-0.3	103	0.00
47		2-Chloronaphthalene	1.075	1.072	0.3	102	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.384	0.397	-3.4	104	0.00
49		Acenaphthylene	1.783	1.790	-0.4	102	0.00
50		Dimethylphthalate	1.479	1.470	0.6	102	0.00
51		2,6-Dinitrotoluene	0.313	0.318	-1.6	103	0.00
52	C	Acenaphthene	1.073	1.073	0.0	103	0.00
53		3-Nitroaniline	0.351	0.362	-3.1	103	0.00
54	P	2,4-Dinitrophenol	0.196	0.202	-3.1	104	0.00
55		Dibenzofuran	1.732	1.716	0.9	103	0.00
56	P	4-Nitrophenol	0.278	0.286	-2.9	103	0.00
57		2,4-Dinitrotoluene	0.415	0.430	-3.6	103	0.00
58		Fluorene	1.365	1.351	1.0	103	0.00
59		2,3,4,6-Tetrachlorophenol	0.379	0.382	-0.8	103	0.00
60		Diethylphthalate	1.471	1.466	0.3	102	0.00
61		4-Chlorophenyl-phenylether	0.716	0.717	-0.1	103	0.00
62		4-Nitroaniline	0.337	0.347	-3.0	102	0.00
63		Azobenzene	1.491	1.493	-0.1	102	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
65		4,6-Dinitro-2-methylphenol	0.115	0.123	-7.0	104	0.00
66	c	n-Nitrosodiphenylamine	0.565	0.561	0.7	102	0.00
67		4-Bromophenyl-phenylether	0.206	0.206	0.0	102	0.00
68		Hexachlorobenzene	0.218	0.218	0.0	103	0.00
69		Atrazine	0.176	0.170	3.4	103	0.00
70	C	Pentachlorophenol	0.165	0.168	-1.8	103	0.00
71		Phenanthrene	0.997	0.990	0.7	103	0.00
72		Anthracene	1.018	1.016	0.2	102	0.00
73		Carbazole	0.933	0.931	0.2	102	0.00
74		Di-n-butylphthalate	1.198	1.209	-0.9	102	0.00
75	C	Fluoranthene	1.171	1.178	-0.6	103	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
77		Benzidine	0.408	0.404	1.0	108	0.00
78		Pyrene	1.392	1.399	-0.5	103	0.00
79	S	Terphenyl-d14	1.098	1.123	-2.3	102	0.00
80		Butylbenzylphthalate	0.619	0.622	-0.5	101	0.00
81		Benzo(a)anthracene	1.304	1.311	-0.5	103	0.00
82		3,3'-Dichlorobenzidine	0.459	0.460	-0.2	101	0.00
83		Chrysene	1.223	1.211	1.0	102	0.00
84		Bis(2-ethylhexyl)phthalate	0.918	0.929	-1.2	102	0.00
85	c	Di-n-octyl phthalate	1.546	1.557	-0.7	101	0.00
86	I	Perylene-d12	1.000	1.000	0.0	104	0.00
87		Indeno(1,2,3-cd)pyrene	1.194	1.199	-0.4	105	0.00
88		Benzo(b)fluoranthene	1.168	1.159	0.8	101	0.00
89		Benzo(k)fluoranthene	1.154	1.165	-1.0	106	0.00
90	C	Benzo(a)pyrene	1.096	1.100	-0.4	104	0.00
91		Dibenzo(a,h)anthracene	0.996	1.009	-1.3	105	0.00
92		Benzo(g,h,i)perylene	0.951	0.948	0.3	104	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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