Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100421\
Data File : BM032317.D
Acq On : 04 Oct 2021 09:52
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Oct 04 11:02:41 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM100221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 01 18:34:58 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	102	0.00
2		1,4-Dioxane	0.505	0.487	3.6	104	0.00
3		Pyridine	1.367	1.323	3.2	102	0.00
4		n-Nitrosodimethylamine	0.548	0.530	3.3	101	0.00
5	S	2-Fluorophenol	1.187	1.138	4.1	101	0.00
6		Aniline	1.794	1.755	2.2	101	0.00
7	S	Phenol-d6	1.623	1.582	2.5	101	0.00
8		2-Chlorophenol	1.294	1.253	3.2	101	0.00
9		Benzaldehyde	0.947	0.857	9.5	102	0.00
10	C	Phenol	1.563	1.527	2.3	102	0.00
11		bis(2-Chloroethyl)ether	1.243	1.220	1.9	103	0.00
12		1,3-Dichlorobenzene	1.458	1.397	4.2	101	0.00
13	C	1,4-Dichlorobenzene	1.484	1.421	4.2	101	0.00
14		1,2-Dichlorobenzene	1.425	1.372	3.7	102	0.00
15		Benzyl Alcohol	1.112	1.119	-0.6	103	0.00
16		2,2'-oxybis(1-Chloropropane	1.575	1.558	1.1	103	0.00
17		2-Methylphenol	1.056	1.055	0.1	103	0.00
18		Hexachloroethane	0.505	0.504	0.2	102	0.00
19	P	n-Nitroso-di-n-propylamine	0.883	0.883	0.0	102	0.00
20		3+4-Methylphenols	1.425	1.429	-0.3	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.482	0.466	3.3	103	0.00
23	S	Nitrobenzene-d5	0.360	0.353	1.9	102	0.00
24		Nitrobenzene	0.347	0.340	2.0	103	0.00
25		Isophorone	0.598	0.601	-0.5	103	0.00
26	C	2-Nitrophenol	0.161	0.165	-2.5	103	0.00
27		2,4-Dimethylphenol	0.272	0.268	1.5	103	0.00
28		bis(2-Chloroethoxy)methane	0.430	0.422	1.9	104	0.00
29	C	2,4-Dichlorophenol	0.303	0.301	0.7	103	0.00
30		1,2,4-Trichlorobenzene	0.344	0.328	4.7	102	0.00
31		Naphthalene	1.026	0.987	3.8	103	0.00
32		Benzoic acid	0.189	0.195	-3.2	99	0.00
33		4-Chloroaniline	0.424	0.416	1.9	102	0.00
34	C	Hexachlorobutadiene	0.206	0.196	4.9	102	0.00
35		Caprolactam	0.093	0.094	-1.1	97	0.00
36	C	4-Chloro-3-methylphenol	0.326	0.324	0.6	102	0.00
37		2-Methylnaphthalene	0.697	0.677	2.9	102	0.00
38		1-Methylnaphthalene	0.682	0.660	3.2	102	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
40		1,2,4,5-Tetrachlorobenzene	0.576	0.547	5.0	102	0.00
41	P	Hexachlorocyclopentadiene	0.275	0.278	-1.1	103	0.00
42	S	2,4,6-Tribromophenol	0.224	0.214	4.5	98	0.00
43	C	2,4,6-Trichlorophenol	0.399	0.395	1.0	103	0.00
44		2,4,5-Trichlorophenol	0.428	0.420	1.9	101	0.00
45	S	2-Fluorobiphenyl	1.345	1.218	9.4	101	0.00
46		1,1'-Biphenyl	1.476	1.409	4.5	102	0.00
47		2-Chloronaphthalene	1.132	1.087	4.0	102	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.308	0.318	-3.2	100	0.00
49		Acenaphthylene	1.757	1.718	2.2	102	0.00
50		Dimethylphthalate	1.412	1.358	3.8	100	0.00
51		2,6-Dinitrotoluene	0.286	0.285	0.3	101	0.00
52	C	Acenaphthene	1.159	1.124	3.0	102	0.00
53		3-Nitroaniline	0.321	0.321	0.0	99	0.00
54	P	2,4-Dinitrophenol	0.159	0.166	-4.4	102	0.00
55		Dibenzofuran	1.765	1.676	5.0	100	0.00
56	P	4-Nitrophenol	0.265	0.264	0.4	97	0.00
57		2,4-Dinitrotoluene	0.377	0.384	-1.9	99	0.00
58		Fluorene	1.342	1.212	9.7	101	0.00
59		2,3,4,6-Tetrachlorophenol	0.369	0.359	2.7	99	0.00
60		Diethylphthalate	1.387	1.343	3.2	101	0.00
61		4-Chlorophenyl-phenylether	0.695	0.626	9.9	101	0.00
62		4-Nitroaniline	0.321	0.319	0.6	97	0.00
63		Azobenzene	1.320	1.305	1.1	101	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
65		4,6-Dinitro-2-methylphenol	0.113	0.121	-7.1	99	0.00
66	c	n-Nitrosodiphenylamine	0.591	0.576	2.5	99	0.00
67		4-Bromophenyl-phenylether	0.208	0.203	2.4	99	0.00
68		Hexachlorobenzene	0.229	0.220	3.9	98	0.00
69		Atrazine	0.201	0.196	2.5	98	0.00
70	C	Pentachlorophenol	0.141	0.143	-1.4	97	0.00
71		Phenanthrene	1.050	0.998	5.0	98	0.00
72		Anthracene	1.026	0.993	3.2	98	0.00
73		Carbazole	1.019	0.976	4.2	96	0.00
74		Di-n-butylphthalate	1.094	1.103	-0.8	96	0.00
75	C	Fluoranthene	1.164	1.094	6.0	94	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
77		Benzidine	0.497	0.463	6.8	83	0.00
78		Pyrene	1.356	1.352	0.3	94	0.00
79	S	Terphenyl-d14	0.950	0.910	4.2	94	0.00
80		Butylbenzylphthalate	0.515	0.553	-7.4	91	0.00
81		Benzo(a)anthracene	1.250	1.206	3.5	90	0.00
82		3,3'-Dichlorobenzidine	0.394	0.390	1.0	88	0.00
83		Chrysene	1.209	1.157	4.3	89	0.00
84		Bis(2-ethylhexyl)phthalate	0.695	0.732	-5.3	90	0.00
85	c	Di-n-octyl phthalate	1.165	1.216	-4.4	85	0.00
86	I	Perylene-d12	1.000	1.000	0.0	83	0.00
87		Indeno(1,2,3-cd)pyrene	1.232	1.167	5.3	81	0.00
88		Benzo(b)fluoranthene	1.217	1.170	3.9	82	0.00
89		Benzo(k)fluoranthene	1.189	1.171	1.5	87	0.00
90	C	Benzo(a)pyrene	0.987	0.974	1.3	84	0.00
91		Dibenzo(a,h)anthracene	1.039	0.979	5.8	81	-0.01
92		Benzo(g,h,i)perylene	1.027	0.972	5.4	82	0.00

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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