Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032923\
Data File : BM039208.D
Acq On : 29 Mar 2023 16:37
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM032923

Quant Time: Mar 30 05:17:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 30 05:15:45 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	87	0.00
2		1,4-Dioxane	0.543	0.536	1.3	87	0.00
3		Pyridine	1.520	1.505	1.0	84	0.00
4		n-Nitrosodimethylamine	0.597	0.588	1.5	86	0.00
5	S	2-Fluorophenol	1.327	1.323	0.3	87	0.00
6		Aniline	2.125	2.108	0.8	86	0.00
7	S	Phenol-d6	1.717	1.731	-0.8	87	0.00
8		2-Chlorophenol	1.359	1.361	-0.1	87	0.00
9		Benzaldehyde	0.978	0.995	-1.7	89	0.00
10	C	Phenol	1.721	1.732	-0.6	88	0.00
11		bis(2-Chloroethyl)ether	1.384	1.376	0.6	87	0.00
12		1,3-Dichlorobenzene	1.441	1.425	1.1	86	0.00
13	C	1,4-Dichlorobenzene	1.457	1.437	1.4	86	0.00
14		1,2-Dichlorobenzene	1.393	1.382	0.8	87	0.00
15		Benzyl Alcohol	1.181	1.191	-0.8	87	0.00
16		2,2'-oxybis(1-Chloropropane	1.771	1.757	0.8	87	-0.01
17		2-Methylphenol	1.204	1.200	0.3	87	0.00
18		Hexachloroethane	0.564	0.557	1.2	87	0.00
19	P	n-Nitroso-di-n-propylamine	1.098	1.091	0.6	87	0.00
20		3+4-Methylphenols	1.618	1.634	-1.0	87	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	87	0.00
22		Acetophenone	0.535	0.527	1.5	85	0.00
23	S	Nitrobenzene-d5	0.412	0.411	0.2	86	0.00
24		Nitrobenzene	0.409	0.406	0.7	86	0.00
25		Isophorone	0.768	0.770	-0.3	86	0.00
26	C	2-Nitrophenol	0.189	0.191	-1.1	86	0.00
27		2,4-Dimethylphenol	0.314	0.318	-1.3	87	0.00
28		bis(2-Chloroethoxy)methane	0.465	0.464	0.2	87	0.00
29	C	2,4-Dichlorophenol	0.307	0.314	-2.3	87	0.00
30		1,2,4-Trichlorobenzene	0.328	0.327	0.3	87	0.00
31		Naphthalene	1.078	1.080	-0.2	87	0.00
32		Benzoic acid	0.245	0.244	0.4	83	0.00
33		4-Chloroaniline	0.472	0.476	-0.8	87	0.00
34	C	Hexachlorobutadiene	0.193	0.193	0.0	87	0.00
35		Caprolactam	0.106	0.108	-1.9	85	0.00
36	C	4-Chloro-3-methylphenol	0.338	0.343	-1.5	86	0.00
37		2-Methylnaphthalene	0.750	0.753	-0.4	87	0.00
38		1-Methylnaphthalene	0.702	0.699	0.4	86	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	88	0.00
40		1,2,4,5-Tetrachlorobenzene	0.609	0.612	-0.5	87	0.00
41	P	Hexachlorocyclopentadiene	0.333	0.355	-6.6	88	0.00
42	S	2,4,6-Tribromophenol	0.254	0.257	-1.2	87	0.00
43	C	2,4,6-Trichlorophenol	0.415	0.426	-2.7	87	0.00
44		2,4,5-Trichlorophenol	0.487	0.500	-2.7	87	0.00
45	S	2-Fluorobiphenyl	1.476	1.482	-0.4	88	0.00
46		1,1'-Biphenyl	1.579	1.584	-0.3	87	0.00
47		2-Chloronaphthalene	1.196	1.195	0.1	87	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.402	0.408	-1.5	87	0.00
49		Acenaphthylene	1.965	1.962	0.2	87	0.00
50		Dimethylphthalate	1.541	1.544	-0.2	88	0.00
51		2,6-Dinitrotoluene	0.331	0.340	-2.7	88	0.00
52	C	Acenaphthene	1.161	1.159	0.2	87	0.00
53		3-Nitroaniline	0.383	0.390	-1.8	86	0.00
54	P	2,4-Dinitrophenol	0.196	0.200	-2.0	86	0.00
55		Dibenzofuran	1.868	1.867	0.1	88	0.00
56	P	4-Nitrophenol	0.295	0.306	-3.7	84	0.00
57		2,4-Dinitrotoluene	0.448	0.464	-3.6	87	0.00
58		Fluorene	1.480	1.489	-0.6	88	0.00
59		2,3,4,6-Tetrachlorophenol	0.387	0.390	-0.8	87	0.00
60		Diethylphthalate	1.551	1.558	-0.5	88	0.00
61		4-Chlorophenyl-phenylether	0.733	0.736	-0.4	88	0.00
62		4-Nitroaniline	0.393	0.403	-2.5	85	0.00
63		Azobenzene	1.554	1.570	-1.0	88	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	88	0.00
65		4,6-Dinitro-2-methylphenol	0.132	0.135	-2.3	87	0.00
66	c	n-Nitrosodiphenylamine	0.630	0.629	0.2	87	0.00
67		4-Bromophenyl-phenylether	0.214	0.216	-0.9	88	0.00
68		Hexachlorobenzene	0.232	0.233	-0.4	88	0.00
69		Atrazine	0.200	0.202	-1.0	88	0.00
70	C	Pentachlorophenol	0.166	0.174	-4.8	87	0.00
71		Phenanthrene	1.095	1.105	-0.9	89	0.00
72		Anthracene	1.117	1.118	-0.1	87	0.00
73		Carbazole	1.044	1.044	0.0	86	0.00
74		Di-n-butylphthalate	1.291	1.326	-2.7	88	0.00
75	C	Fluoranthene	1.249	1.254	-0.4	87	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	85	0.00
77		Benzidine	0.528	0.542	-2.7	82	0.00
78		Pyrene	1.424	1.460	-2.5	87	0.00
79	S	Terphenyl-d14	1.083	1.119	-3.3	88	0.00
80		Butylbenzylphthalate	0.652	0.670	-2.8	86	0.00
81		Benzo(a)anthracene	1.411	1.412	-0.1	86	0.00
82		3,3'-Dichlorobenzidine	0.470	0.471	-0.2	85	0.00
83		Chrysene	1.362	1.373	-0.8	85	0.00
84		Bis(2-ethylhexyl)phthalate	0.963	0.995	-3.3	88	0.00
85	c	Di-n-octyl phthalate	1.736	1.774	-2.2	87	0.00
86	I	Perylene-d12	1.000	1.000	0.0	85	0.00
87		Indeno(1,2,3-cd)pyrene	1.474	1.481	-0.5	84	0.00
88		Benzo(b)fluoranthene	1.224	1.237	-1.1	86	0.00
89		Benzo(k)fluoranthene	1.247	1.267	-1.6	86	0.00
90	C	Benzo(a)pyrene	1.202	1.216	-1.2	85	0.00
91		Dibenzo(a,h)anthracene	1.237	1.243	-0.5	84	0.00
92		Benzo(g,h,i)perylene	1.217	1.218	-0.1	83	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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