Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081024\
Data File : BM047148.D
Acq On : 10 Aug 2024 19:10
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM081024

Quant Time: Aug 11 23:52:13 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Aug 11 23:47:58 2024
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	86	0.00
2		1,4-Dioxane	0.464	0.427	8.0	84	0.00
3		Pyridine	1.360	1.291	5.1	86	0.00
4		n-Nitrosodimethylamine	0.594	0.554	6.7	84	0.00
5	S	2-Fluorophenol	1.119	1.052	6.0	85	0.00
6		Aniline	1.458	1.425	2.3	86	0.00
7	S	Phenol-d6	1.542	1.477	4.2	86	0.00
8		2-Chlorophenol	1.221	1.188	2.7	87	0.00
9		Benzaldehyde	0.904	0.804	11.1	89	0.00
10	C	Phenol	1.539	1.445	6.1	86	0.00
11		bis(2-Chloroethyl)ether	1.318	1.251	5.1	84	0.00
12		1,3-Dichlorobenzene	1.451	1.411	2.8	87	0.00
13	C	1,4-Dichlorobenzene	1.469	1.426	2.9	86	0.00
14		1,2-Dichlorobenzene	1.386	1.342	3.2	88	0.00
15		Benzyl Alcohol	1.098	1.086	1.1	85	0.00
16		2,2'-oxybis(1-Chloropropane	1.691	1.598	5.5	83	0.00
17		2-Methylphenol	1.027	1.005	2.1	87	0.00
18		Hexachloroethane	0.569	0.547	3.9	86	0.00
19	P	n-Nitroso-di-n-propylamine	1.039	0.983	5.4	86	0.00
20		3+4-Methylphenols	1.435	1.425	0.7	87	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	87	0.00
22		Acetophenone	0.483	0.460	4.8	86	0.00
23	S	Nitrobenzene-d5	0.397	0.383	3.5	85	0.00
24		Nitrobenzene	0.400	0.390	2.5	86	0.00
25		Isophorone	0.692	0.684	1.2	86	0.00
26	C	2-Nitrophenol	0.177	0.183	-3.4	87	0.00
27		2,4-Dimethylphenol	0.207	0.205	1.0	85	0.00
28		bis(2-Chloroethoxy)methane	0.435	0.425	2.3	85	0.00
29	C	2,4-Dichlorophenol	0.308	0.315	-2.3	88	0.00
30		1,2,4-Trichlorobenzene	0.348	0.351	-0.9	88	0.00
31		Naphthalene	0.996	0.961	3.5	86	0.00
32		Benzoic acid	0.241	0.260	-7.9	90	-0.01
33		4-Chloroaniline	0.383	0.386	-0.8	87	0.00
34	C	Hexachlorobutadiene	0.204	0.205	-0.5	88	0.00
35		Caprolactam	0.107	0.108	-0.9	86	0.00
36	C	4-Chloro-3-methylphenol	0.328	0.326	0.6	87	0.00
37		2-Methylnaphthalene	0.709	0.691	2.5	87	0.00
38		1-Methylnaphthalene	0.701	0.681	2.9	87	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	87	0.00
40		1,2,4,5-Tetrachlorobenzene	0.546	0.553	-1.3	89	0.00
41	P	Hexachlorocyclopentadiene	0.190	0.195	-2.6	85	0.00
42	S	2,4,6-Tribromophenol	0.232	0.237	-2.2	89	0.00
43	C	2,4,6-Trichlorophenol	0.388	0.395	-1.8	87	0.00
44		2,4,5-Trichlorophenol	0.430	0.442	-2.8	88	0.00
45	S	2-Fluorobiphenyl	1.297	1.276	1.6	88	0.00
46		1,1'-Biphenyl	1.424	1.375	3.4	87	0.00
47		2-Chloronaphthalene	1.113	1.070	3.9	87	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.349	0.351	-0.6	86	0.00
49		Acenaphthylene	1.640	1.610	1.8	88	0.00
50		Dimethylphthalate	1.397	1.375	1.6	87	0.00
51		2,6-Dinitrotoluene	0.327	0.329	-0.6	88	0.00
52	C	Acenaphthene	1.041	1.010	3.0	88	0.00
53		3-Nitroaniline	0.327	0.331	-1.2	87	0.00
54	P	2,4-Dinitrophenol	0.196	0.197	-0.5	85	0.00
55		Dibenzofuran	1.644	1.594	3.0	88	0.00
56	P	4-Nitrophenol	0.282	0.292	-3.5	87	0.00
57		2,4-Dinitrotoluene	0.436	0.443	-1.6	88	0.00
58		Fluorene	1.356	1.317	2.9	88	0.00
59		2,3,4,6-Tetrachlorophenol	0.354	0.361	-2.0	89	0.00
60		Diethylphthalate	1.426	1.395	2.2	87	0.00
61		4-Chlorophenyl-phenylether	0.646	0.626	3.1	88	0.00
62		4-Nitroaniline	0.322	0.323	-0.3	89	0.00
63		Azobenzene	1.352	1.280	5.3	85	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	88	0.00
65		4,6-Dinitro-2-methylphenol	0.117	0.124	-6.0	87	0.00
66	c	n-Nitrosodiphenylamine	0.545	0.535	1.8	88	0.00
67		4-Bromophenyl-phenylether	0.193	0.193	0.0	89	0.00
68		Hexachlorobenzene	0.231	0.232	-0.4	89	0.00
69		Atrazine	0.164	0.169	-3.0	87	0.00
70	C	Pentachlorophenol	0.162	0.168	-3.7	88	0.00
71		Phenanthrene	1.004	0.978	2.6	88	0.00
72		Anthracene	0.977	0.967	1.0	89	0.00
73		Carbazole	0.997	0.975	2.2	88	0.00
74		Di-n-butylphthalate	1.160	1.151	0.8	87	0.00
75	C	Fluoranthene	1.221	1.190	2.5	87	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	88	0.00
77		Benzidine	0.339	0.503	-48.4#	91	0.00
78		Pyrene	1.428	1.404	1.7	87	0.00
79	S	Terphenyl-d14	0.933	0.930	0.3	88	0.00
80		Butylbenzylphthalate	0.583	0.602	-3.3	87	0.00
81		Benzo(a)anthracene	1.328	1.297	2.3	88	0.00
82		3,3'-Dichlorobenzidine	0.415	0.447	-7.7	89	0.00
83		Chrysene	1.260	1.217	3.4	88	0.00
84		Bis(2-ethylhexyl)phthalate	0.826	0.834	-1.0	87	0.00
85	c	Di-n-octyl phthalate	1.405	1.451	-3.3	88	0.00
86	I	Perylene-d12	1.000	1.000	0.0	90	0.00
87		Indeno(1,2,3-cd)pyrene	1.293	1.310	-1.3	92	0.00
88		Benzo(b)fluoranthene	1.188	1.141	4.0	88	0.00
89		Benzo(k)fluoranthene	1.122	1.104	1.6	90	0.00
90	C	Benzo(a)pyrene	1.029	1.023	0.6	90	0.00
91		Dibenzo(a,h)anthracene	1.046	1.049	-0.3	92	0.00
92		Benzo(g,h,i)perylene	1.086	1.103	-1.6	92	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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