Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102622\
Data File : BM037241.D
Acq On : 26 Oct 2022 16:36
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM102622

Quant Time: Oct 27 04:58:59 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 27 04:49:27 2022
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	113	0.00
2		1,4-Dioxane	0.400	0.386	3.5	108	0.00
3		Pyridine	1.171	1.177	-0.5	107	0.00
4		n-Nitrosodimethylamine	0.490	0.471	3.9	105	0.00
5	S	2-Fluorophenol	0.976	0.988	-1.2	111	0.00
6		Aniline	1.682	1.751	-4.1	110	0.00
7	S	Phenol-d6	1.350	1.376	-1.9	110	0.00
8		2-Chlorophenol	1.290	1.352	-4.8	113	0.00
9		Benzaldehyde	0.745	0.723	3.0	113	0.00
10	C	Phenol	1.504	1.531	-1.8	109	0.00
11		bis(2-Chloroethyl)ether	1.222	1.226	-0.3	108	0.00
12		1,3-Dichlorobenzene	1.443	1.481	-2.6	112	0.00
13	C	1,4-Dichlorobenzene	1.490	1.544	-3.6	113	0.00
14		1,2-Dichlorobenzene	1.441	1.491	-3.5	113	0.00
15		Benzyl Alcohol	1.020	1.095	-7.4	112	0.00
16		2,2'-oxybis(1-Chloropropane	1.605	1.602	0.2	109	0.00
17		2-Methylphenol	1.049	1.110	-5.8	113	0.00
18		Hexachloroethane	0.516	0.537	-4.1	112	0.00
19	P	n-Nitroso-di-n-propylamine	0.977	1.025	-4.9	110	0.00
20		3+4-Methylphenols	1.448	1.539	-6.3	113	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.468	0.494	-5.6	113	0.00
23	S	Nitrobenzene-d5	0.350	0.368	-5.1	109	0.00
24		Nitrobenzene	0.358	0.372	-3.9	110	0.00
25		Isophorone	0.609	0.654	-7.4	112	0.00
26	C	2-Nitrophenol	0.164	0.184	-12.2	115	0.00
27		2,4-Dimethylphenol	0.250	0.269	-7.6	113	0.00
28		bis(2-Chloroethoxy)methane	0.367	0.383	-4.4	112	0.00
29	C	2,4-Dichlorophenol	0.307	0.332	-8.1	113	0.00
30		1,2,4-Trichlorobenzene	0.343	0.358	-4.4	113	0.00
31		Naphthalene	1.068	1.107	-3.7	112	0.00
32		Benzoic acid	0.200	0.232	-16.0	121	0.00
33		4-Chloroaniline	0.419	0.453	-8.1	113	0.00
34	C	Hexachlorobutadiene	0.216	0.226	-4.6	113	0.00
35		Caprolactam	0.091	0.102	-12.1	116	0.00
36	C	4-Chloro-3-methylphenol	0.298	0.324	-8.7	113	0.00
37		2-Methylnaphthalene	0.733	0.774	-5.6	113	0.00
38		1-Methylnaphthalene	0.720	0.762	-5.8	113	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	113	0.00
40		1,2,4,5-Tetrachlorobenzene	0.615	0.646	-5.0	114	0.00
41	P	Hexachlorocyclopentadiene	0.309	0.316	-2.3	109	0.00
42	S	2,4,6-Tribromophenol	0.248	0.287	-15.7	116	0.00
43	C	2,4,6-Trichlorophenol	0.408	0.446	-9.3	114	0.00
44		2,4,5-Trichlorophenol	0.452	0.494	-9.3	115	0.00
45	S	2-Fluorobiphenyl	1.416	1.476	-4.2	112	0.00
46		1,1'-Biphenyl	1.472	1.534	-4.2	113	0.00
47		2-Chloronaphthalene	1.168	1.225	-4.9	113	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.307	0.340	-10.7	112	0.00
49		Acenaphthylene	1.832	1.958	-6.9	113	0.00
50		Dimethylphthalate	1.459	1.559	-6.9	113	0.00
51		2,6-Dinitrotoluene	0.303	0.335	-10.6	114	0.00
52	C	Acenaphthene	1.128	1.182	-4.8	112	0.00
53		3-Nitroaniline	0.327	0.373	-14.1	115	0.00
54	P	2,4-Dinitrophenol	0.182	0.207	-13.7	118	0.00
55		Dibenzofuran	1.807	1.888	-4.5	112	0.00
56	P	4-Nitrophenol	0.270	0.294	-8.9	113	0.00
57		2,4-Dinitrotoluene	0.405	0.456	-12.6	114	0.00
58		Fluorene	1.473	1.575	-6.9	112	0.00
59		2,3,4,6-Tetrachlorophenol	0.409	0.448	-9.5	113	0.00
60		Diethylphthalate	1.427	1.532	-7.4	113	0.00
61		4-Chlorophenyl-phenylether	0.733	0.787	-7.4	113	0.00
62		4-Nitroaniline	0.337	0.390	-15.7	115	0.00
63		Azobenzene	1.282	1.363	-6.3	110	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	113	0.00
65		4,6-Dinitro-2-methylphenol	0.123	0.135	-9.8	116	0.00
66	c	n-Nitrosodiphenylamine	0.591	0.627	-6.1	113	0.00
67		4-Bromophenyl-phenylether	0.214	0.229	-7.0	113	0.00
68		Hexachlorobenzene	0.260	0.276	-6.2	114	0.00
69		Atrazine	0.174	0.201	-15.5	114	0.00
70	C	Pentachlorophenol	0.155	0.177	-14.2	116	0.00
71		Phenanthrene	1.114	1.180	-5.9	114	0.00
72		Anthracene	1.064	1.139	-7.0	113	0.00
73		Carbazole	0.970	1.039	-7.1	112	0.00
74		Di-n-butylphthalate	1.127	1.238	-9.8	114	0.00
75	C	Fluoranthene	1.286	1.387	-7.9	113	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	109	0.00
77		Benzidine	0.296	0.344	-16.2	109	0.00
78		Pyrene	1.321	1.423	-7.7	111	0.00
79	S	Terphenyl-d14	0.978	1.073	-9.7	110	0.00
80		Butylbenzylphthalate	0.476	0.533	-12.0	112	0.00
81		Benzo(a)anthracene	1.293	1.368	-5.8	109	0.00
82		3,3'-Dichlorobenzidine	0.380	0.417	-9.7	108	0.00
83		Chrysene	1.240	1.291	-4.1	107	0.00
84		Bis(2-ethylhexyl)phthalate	0.685	0.776	-13.3	111	0.00
85	c	Di-n-octyl phthalate	1.103	1.220	-10.6	109	0.00
86	I	Perylene-d12	1.000	1.000	0.0	104	0.00
87		Indeno(1,2,3-cd)pyrene	1.319	1.411	-7.0	107	0.00
88		Benzo(b)fluoranthene	1.278	1.398	-9.4	106	0.00
89		Benzo(k)fluoranthene	1.318	1.400	-6.2	102	0.00
90	C	Benzo(a)pyrene	1.032	1.101	-6.7	104	0.00
91		Dibenzo(a,h)anthracene	1.110	1.193	-7.5	106	0.00
92		Benzo(g,h,i)perylene	1.044	1.108	-6.1	109	0.00

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