Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060222\
Data File : BM035364.D
Acq On : 02 Jun 2022 16:01
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM060222

Quant Time: Jun 03 00:33:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 02 16:16:23 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	87	0.00
2		1,4-Dioxane	0.319	0.260	18.5	78	0.00
3		Pyridine	0.871	0.746	14.4	78	0.00
4		n-Nitrosodimethylamine	0.329	0.278	15.5	77	0.00
5	S	2-Fluorophenol	0.952	0.858	9.9	86	0.00
6		Aniline	1.421	1.454	-2.3	95	0.00
7	S	Phenol-d6	1.300	1.303	-0.2	95	0.00
8		2-Chlorophenol	1.194	1.197	-0.3	90	0.00
9		Benzaldehyde	0.609	0.579	4.9	99	0.00
10	C	Phenol	1.253	1.259	-0.5	95	0.00
11		bis(2-Chloroethyl)ether	1.020	1.031	-1.1	97	0.00
12		1,3-Dichlorobenzene	1.395	1.373	1.6	87	0.00
13	C	1,4-Dichlorobenzene	1.432	1.412	1.4	87	0.00
14		1,2-Dichlorobenzene	1.404	1.352	3.7	84	0.00
15		Benzyl Alcohol	0.950	0.934	1.7	82	0.00
16		2,2'-oxybis(1-Chloropropane	1.243	1.187	4.5	76	0.00
17		2-Methylphenol	0.948	0.939	0.9	84	0.00
18		Hexachloroethane	0.467	0.418	10.5	77	0.00
19	P	n-Nitroso-di-n-propylamine	0.822	0.750	8.8	75	0.00
20		3+4-Methylphenols	1.326	1.218	8.1	78	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	78	0.00
22		Acetophenone	0.445	0.458	-2.9	78	0.00
23	S	Nitrobenzene-d5	0.315	0.318	-1.0	75	0.00
24		Nitrobenzene	0.285	0.290	-1.8	76	0.00
25		Isophorone	0.545	0.560	-2.8	76	0.00
26	C	2-Nitrophenol	0.187	0.198	-5.9	80	0.00
27		2,4-Dimethylphenol	0.248	0.258	-4.0	79	0.00
28		bis(2-Chloroethoxy)methane	0.376	0.383	-1.9	77	0.00
29	C	2,4-Dichlorophenol	0.338	0.347	-2.7	78	0.00
30		1,2,4-Trichlorobenzene	0.388	0.397	-2.3	79	0.00
31		Naphthalene	0.969	0.984	-1.5	79	0.00
32		Benzoic acid	0.221	0.247	-11.8	78	0.00
33		4-Chloroaniline	0.406	0.424	-4.4	78	0.00
34	C	Hexachlorobutadiene	0.257	0.260	-1.2	79	0.00
35		Caprolactam	0.099	0.105	-6.1	76	0.00
36	C	4-Chloro-3-methylphenol	0.311	0.329	-5.8	78	0.00
37		2-Methylnaphthalene	0.731	0.745	-1.9	77	0.00
38		1-Methylnaphthalene	0.717	0.730	-1.8	77	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	74	0.00
40		1,2,4,5-Tetrachlorobenzene	0.656	0.685	-4.4	78	0.00
41	P	Hexachlorocyclopentadiene	0.281	0.294	-4.6	77	0.00
42	S	2,4,6-Tribromophenol	0.321	0.356	-10.9	85	0.00
43	C	2,4,6-Trichlorophenol	0.443	0.470	-6.1	77	0.00
44		2,4,5-Trichlorophenol	0.475	0.506	-6.5	77	0.00
45	S	2-Fluorobiphenyl	1.392	1.442	-3.6	77	0.00
46		1,1'-Biphenyl	1.426	1.476	-3.5	77	0.00
47		2-Chloronaphthalene	1.112	1.152	-3.6	76	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.247	0.261	-5.7	73	0.00
49		Acenaphthylene	1.694	1.691	0.2	73	0.00
50		Dimethylphthalate	1.470	1.520	-3.4	76	0.00
51		2,6-Dinitrotoluene	0.314	0.322	-2.5	74	0.00
52	C	Acenaphthene	1.026	1.041	-1.5	74	0.00
53		3-Nitroaniline	0.298	0.315	-5.7	74	0.00
54	P	2,4-Dinitrophenol	0.206	0.225	-9.2	77	0.00
55		Dibenzofuran	1.762	1.851	-5.1	78	0.00
56	P	4-Nitrophenol	0.224	0.252	-12.5	76	0.00
57		2,4-Dinitrotoluene	0.430	0.471	-9.5	79	0.00
58		Fluorene	1.376	1.458	-6.0	77	0.00
59		2,3,4,6-Tetrachlorophenol	0.437	0.483	-10.5	81	0.00
60		Diethylphthalate	1.454	1.546	-6.3	77	0.00
61		4-Chlorophenyl-phenylether	0.789	0.824	-4.4	77	0.00
62		4-Nitroaniline	0.311	0.355	-14.1	79	0.00
63		Azobenzene	1.032	1.094	-6.0	75	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	73	0.00
65		4,6-Dinitro-2-methylphenol	0.134	0.158	-17.9	80	0.00
66	c	n-Nitrosodiphenylamine	0.564	0.609	-8.0	78	0.00
67		4-Bromophenyl-phenylether	0.242	0.256	-5.8	78	0.00
68		Hexachlorobenzene	0.271	0.295	-8.9	82	0.00
69		Atrazine	0.208	0.232	-11.5	79	0.00
70	C	Pentachlorophenol	0.148	0.175	-18.2	83	0.00
71		Phenanthrene	1.012	1.099	-8.6	78	0.00
72		Anthracene	1.009	1.092	-8.2	78	0.00
73		Carbazole	0.950	1.045	-10.0	78	0.00
74		Di-n-butylphthalate	1.179	1.304	-10.6	80	0.00
75	C	Fluoranthene	1.222	1.336	-9.3	78	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	82	0.00
77		Benzidine	0.376	0.351	6.6	73	0.00
78		Pyrene	1.213	1.223	-0.8	78	0.00
79	S	Terphenyl-d14	1.004	1.022	-1.8	81	0.00
80		Butylbenzylphthalate	0.515	0.523	-1.6	80	0.00
81		Benzo(a)anthracene	1.228	1.242	-1.1	81	0.00
82		3,3'-Dichlorobenzidine	0.323	0.315	2.5	79	0.00
83		Chrysene	1.172	1.190	-1.5	82	0.00
84		Bis(2-ethylhexyl)phthalate	0.776	0.795	-2.4	81	0.00
85	c	Di-n-octyl phthalate	1.356	1.349	0.5	79	0.00
86	I	Perylene-d12	1.000	1.000	0.0	78	0.00
87		Indeno(1,2,3-cd)pyrene	1.327	1.377	-3.8	84	-0.01
88		Benzo(b)fluoranthene	1.152	1.252	-8.7	80	0.00
89		Benzo(k)fluoranthene	1.179	1.208	-2.5	78	0.00
90	C	Benzo(a)pyrene	0.981	1.016	-3.6	79	0.00
91		Dibenzo(a,h)anthracene	1.107	1.150	-3.9	84	0.00
92		Benzo(g,h,i)perylene	1.068	1.131	-5.9	86	0.00

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