Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071921\
Data File : BM031011.D
Acq On : 19 Jul 2021 18:37
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM071921

Quant Time: Jul 19 19:10:32 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 19 18:48:56 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	106	0.00
2		1,4-Dioxane	0.446	0.434	2.7	104	0.00
3		Pyridine	1.277	1.291	-1.1	105	0.00
4		n-Nitrosodimethylamine	0.723	0.751	-3.9	106	0.00
5	S	2-Fluorophenol	1.122	1.160	-3.4	107	0.00
6		Aniline	1.735	1.818	-4.8	108	0.00
7	S	Phenol-d6	1.614	1.723	-6.8	110	0.00
8		2-Chlorophenol	1.323	1.393	-5.3	108	0.00
9		Benzaldehyde	0.951	1.003	-5.5	112	0.00
10	C	Phenol	1.571	1.650	-5.0	108	0.00
11		bis(2-Chloroethyl)ether	1.169	1.229	-5.1	109	0.00
12		1,3-Dichlorobenzene	1.427	1.461	-2.4	107	0.00
13	C	1,4-Dichlorobenzene	1.457	1.495	-2.6	108	0.00
14		1,2-Dichlorobenzene	1.420	1.487	-4.7	110	0.00
15		Benzyl Alcohol	1.370	1.480	-8.0	110	0.00
16		2,2'-oxybis(1-Chloropropane	1.469	1.522	-3.6	107	0.00
17		2-Methylphenol	1.117	1.193	-6.8	109	0.00
18		Hexachloroethane	0.568	0.599	-5.5	106	0.00
19	P	n-Nitroso-di-n-propylamine	1.163	1.243	-6.9	109	0.00
20		3+4-Methylphenols	1.559	1.661	-6.5	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
22		Acetophenone	0.500	0.513	-2.6	108	0.00
23	S	Nitrobenzene-d5	0.399	0.412	-3.3	108	0.00
24		Nitrobenzene	0.404	0.411	-1.7	108	0.00
25		Isophorone	0.720	0.753	-4.6	109	0.00
26	C	2-Nitrophenol	0.172	0.183	-6.4	110	0.00
27		2,4-Dimethylphenol	0.272	0.280	-2.9	109	0.00
28		bis(2-Chloroethoxy)methane	0.367	0.378	-3.0	108	0.00
29	C	2,4-Dichlorophenol	0.319	0.336	-5.3	109	0.00
30		1,2,4-Trichlorobenzene	0.346	0.354	-2.3	108	0.00
31		Naphthalene	1.032	1.066	-3.3	108	0.00
32		Benzoic acid	0.221	0.250	-13.1	116	0.00
33		4-Chloroaniline	0.442	0.462	-4.5	109	0.00
34	C	Hexachlorobutadiene	0.226	0.227	-0.4	106	0.00
35		Caprolactam	0.105	0.113	-7.6	111	0.00
36	C	4-Chloro-3-methylphenol	0.364	0.383	-5.2	108	0.00
37		2-Methylnaphthalene	0.792	0.818	-3.3	108	0.00
38		1-Methylnaphthalene	0.757	0.786	-3.8	109	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
40		1,2,4,5-Tetrachlorobenzene	0.568	0.586	-3.2	108	0.00
41	P	Hexachlorocyclopentadiene	0.321	0.330	-2.8	102	0.00
42	S	2,4,6-Tribromophenol	0.262	0.282	-7.6	110	0.00
43	C	2,4,6-Trichlorophenol	0.397	0.421	-6.0	109	0.00
44		2,4,5-Trichlorophenol	0.435	0.461	-6.0	108	0.00
45	S	2-Fluorobiphenyl	1.359	1.407	-3.5	108	0.00
46		1,1'-Biphenyl	1.424	1.474	-3.5	108	0.00
47		2-Chloronaphthalene	1.129	1.164	-3.1	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.354	0.384	-8.5	109	0.00
49		Acenaphthylene	1.793	1.867	-4.1	108	0.00
50		Dimethylphthalate	1.499	1.550	-3.4	107	0.00
51		2,6-Dinitrotoluene	0.333	0.349	-4.8	106	0.00
52	C	Acenaphthene	1.119	1.158	-3.5	108	0.00
53		3-Nitroaniline	0.337	0.361	-7.1	108	0.00
54	P	2,4-Dinitrophenol	0.194	0.221	-13.9	110	0.00
55		Dibenzofuran	1.839	1.879	-2.2	106	0.00
56	P	4-Nitrophenol	0.296	0.319	-7.8	109	0.00
57		2,4-Dinitrotoluene	0.468	0.504	-7.7	108	0.00
58		Fluorene	1.550	1.610	-3.9	107	0.00
59		2,3,4,6-Tetrachlorophenol	0.406	0.427	-5.2	108	0.00
60		Diethylphthalate	1.610	1.674	-4.0	107	0.00
61		4-Chlorophenyl-phenylether	0.780	0.795	-1.9	106	0.00
62		4-Nitroaniline	0.368	0.397	-7.9	107	0.00
63		Azobenzene	1.590	1.641	-3.2	106	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
65		4,6-Dinitro-2-methylphenol	0.125	0.137	-9.6	108	0.00
66	c	n-Nitrosodiphenylamine	0.597	0.612	-2.5	106	0.00
67		4-Bromophenyl-phenylether	0.211	0.216	-2.4	106	0.00
68		Hexachlorobenzene	0.236	0.245	-3.8	107	0.00
69		Atrazine	0.216	0.225	-4.2	106	0.00
70	C	Pentachlorophenol	0.154	0.168	-9.1	108	0.00
71		Phenanthrene	1.106	1.132	-2.4	106	0.00
72		Anthracene	1.091	1.127	-3.3	106	0.00
73		Carbazole	1.070	1.103	-3.1	106	0.00
74		Di-n-butylphthalate	1.274	1.350	-6.0	106	0.00
75	C	Fluoranthene	1.363	1.406	-3.2	105	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
77		Benzidine	0.520	0.583	-12.1	88	0.00
78		Pyrene	1.273	1.345	-5.7	105	0.00
79	S	Terphenyl-d14	1.027	1.095	-6.6	105	0.00
80		Butylbenzylphthalate	0.545	0.600	-10.1	106	0.00
81		Benzo(a)anthracene	1.323	1.387	-4.8	105	0.00
82		3,3'-Dichlorobenzidine	0.360	0.390	-8.3	105	0.00
83		Chrysene	1.287	1.315	-2.2	103	0.00
84		Bis(2-ethylhexyl)phthalate	0.838	0.919	-9.7	105	0.00
85	c	Di-n-octyl phthalate	1.417	1.530	-8.0	104	0.00
86	I	Perylene-d12	1.000	1.000	0.0	104	0.00
87		Indeno(1,2,3-cd)pyrene	1.391	1.456	-4.7	104	0.00
88		Benzo(b)fluoranthene	1.332	1.377	-3.4	100	0.00
89		Benzo(k)fluoranthene	1.269	1.351	-6.5	106	0.00
90	C	Benzo(a)pyrene	1.202	1.263	-5.1	103	0.00
91		Dibenzo(a,h)anthracene	1.202	1.256	-4.5	104	0.00
92		Benzo(g,h,i)perylene	1.140	1.181	-3.6	104	0.00

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