Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042121\
Data File : BM029570.D
Acq On : 21 Apr 2021 15:03
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM042121

Quant Time: Apr 21 15:54:32 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 21 15:14:18 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	103	0.01
2		1,4-Dioxane	0.402	0.394	2.0	100	0.01
3		Pyridine	1.025	1.085	-5.9	103	0.01
4		n-Nitrosodimethylamine	0.503	0.489	2.8	96	0.01
5	S	2-Fluorophenol	0.998	1.013	-1.5	103	0.02
6		Aniline	1.638	1.718	-4.9	103	0.01
7	S	Phenol-d6	1.490	1.548	-3.9	102	0.02
8		2-Chlorophenol	1.240	1.261	-1.7	102	0.02
9		Benzaldehyde	0.847	0.827	2.4	106	0.02
10	C	Phenol	1.444	1.490	-3.2	102	0.01
11		bis(2-Chloroethyl)ether	1.102	1.136	-3.1	101	0.01
12		1,3-Dichlorobenzene	1.429	1.440	-0.8	102	0.02
13	C	1,4-Dichlorobenzene	1.464	1.464	0.0	103	0.01
14		1,2-Dichlorobenzene	1.459	1.464	-0.3	102	0.02
15		Benzyl Alcohol	1.083	1.149	-6.1	101	0.01
16		2,2'-oxybis(1-Chloropropane	1.452	1.443	0.6	100	0.02
17		2-Methylphenol	1.006	1.053	-4.7	101	0.01
18		Hexachloroethane	0.529	0.535	-1.1	103	0.01
19	P	n-Nitroso-di-n-propylamine	1.087	1.136	-4.5	102	0.02
20		3+4-Methylphenols	1.384	1.462	-5.6	103	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	101	0.01
22		Acetophenone	0.480	0.504	-5.0	105	0.01
23	S	Nitrobenzene-d5	0.388	0.408	-5.2	102	0.02
24		Nitrobenzene	0.393	0.409	-4.1	102	0.02
25		Isophorone	0.730	0.763	-4.5	101	0.01
26	C	2-Nitrophenol	0.170	0.187	-10.0	103	0.01
27		2,4-Dimethylphenol	0.262	0.278	-6.1	102	0.01
28		bis(2-Chloroethoxy)methane	0.361	0.376	-4.2	102	0.01
29	C	2,4-Dichlorophenol	0.346	0.368	-6.4	102	0.01
30		1,2,4-Trichlorobenzene	0.398	0.409	-2.8	102	0.01
31		Naphthalene	1.028	1.042	-1.4	101	0.02
32		Benzoic acid	0.202	0.231	-14.4	106	0.02
33		4-Chloroaniline	0.406	0.436	-7.4	102	0.01
34	C	Hexachlorobutadiene	0.254	0.261	-2.8	103	0.01
35		Caprolactam	0.105	0.112	-6.7	104	0.02
36	C	4-Chloro-3-methylphenol	0.338	0.360	-6.5	103	0.01
37		2-Methylnaphthalene	0.823	0.841	-2.2	101	0.02
38		1-Methylnaphthalene	0.785	0.802	-2.2	102	0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	102	0.01
40		1,2,4,5-Tetrachlorobenzene	0.600	0.631	-5.2	103	0.01
41	P	Hexachlorocyclopentadiene	0.335	0.347	-3.6	102	0.01
42	S	2,4,6-Tribromophenol	0.277	0.293	-5.8	103	0.02
43	C	2,4,6-Trichlorophenol	0.412	0.449	-9.0	103	0.01
44		2,4,5-Trichlorophenol	0.455	0.488	-7.3	102	0.01
45	S	2-Fluorobiphenyl	1.474	1.524	-3.4	101	0.01
46		1,1'-Biphenyl	1.462	1.517	-3.8	102	0.01
47		2-Chloronaphthalene	1.170	1.219	-4.2	101	0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.294	0.330	-12.2	101	0.01
49		Acenaphthylene	1.794	1.885	-5.1	101	0.01
50		Dimethylphthalate	1.512	1.588	-5.0	103	0.01
51		2,6-Dinitrotoluene	0.333	0.358	-7.5	103	0.01
52	C	Acenaphthene	1.124	1.165	-3.6	102	0.01
53		3-Nitroaniline	0.278	0.315	-13.3	104	0.01
54	P	2,4-Dinitrophenol	0.183	0.207	-13.1	102	0.01
55		Dibenzofuran	1.897	1.968	-3.7	102	0.01
56	P	4-Nitrophenol	0.246	0.268	-8.9	105	0.01
57		2,4-Dinitrotoluene	0.449	0.497	-10.7	104	0.01
58		Fluorene	1.593	1.669	-4.8	102	0.02
59		2,3,4,6-Tetrachlorophenol	0.405	0.442	-9.1	105	0.01
60		Diethylphthalate	1.484	1.559	-5.1	102	0.01
61		4-Chlorophenyl-phenylether	0.806	0.836	-3.7	102	0.01
62		4-Nitroaniline	0.291	0.324	-11.3	101	0.01
63		Azobenzene	1.409	1.492	-5.9	102	0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.01
65		4,6-Dinitro-2-methylphenol	0.133	0.140	-5.3	103	0.01
66	c	n-Nitrosodiphenylamine	0.618	0.640	-3.6	101	0.01
67		4-Bromophenyl-phenylether	0.209	0.220	-5.3	102	0.01
68		Hexachlorobenzene	0.237	0.244	-3.0	102	0.01
69		Atrazine	0.221	0.239	-8.1	102	0.01
70	C	Pentachlorophenol	0.155	0.163	-5.2	106	0.01
71		Phenanthrene	1.152	1.179	-2.3	102	0.02
72		Anthracene	1.136	1.181	-4.0	102	0.01
73		Carbazole	1.064	1.115	-4.8	102	0.01
74		Di-n-butylphthalate	1.156	1.227	-6.1	101	0.01
75	C	Fluoranthene	1.396	1.449	-3.8	102	0.01
76	I	Chrysene-d12	1.000	1.000	0.0	102	0.02
77		Benzidine	0.423	0.472	-11.6	98	0.01
78		Pyrene	1.287	1.297	-0.8	102	0.01
79	S	Terphenyl-d14	1.055	1.058	-0.3	101	0.01
80		Butylbenzylphthalate	0.482	0.503	-4.4	102	0.01
81		Benzo(a)anthracene	1.323	1.324	-0.1	102	0.01
82		3,3'-Dichlorobenzidine	0.423	0.431	-1.9	101	0.02
83		Chrysene	1.300	1.293	0.5	101	0.01
84		Bis(2-ethylhexyl)phthalate	0.772	0.804	-4.1	101	0.01
85	c	Di-n-octyl phthalate	1.295	1.342	-3.6	102	0.01
86	I	Perylene-d12	1.000	1.000	0.0	103	0.02
87		Indeno(1,2,3-cd)pyrene	1.546	1.595	-3.2	103	0.03
88		Benzo(b)fluoranthene	1.304	1.328	-1.8	102	0.02
89		Benzo(k)fluoranthene	1.307	1.325	-1.4	103	0.02
90	C	Benzo(a)pyrene	1.225	1.237	-1.0	102	0.02
91		Dibenzo(a,h)anthracene	1.333	1.380	-3.5	104	0.03
92		Benzo(g,h,i)perylene	1.233	1.275	-3.4	103	0.04

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