Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040921\
Data File : BP005257.D
Acq On : 09 Apr 2021 18:10
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP040921

Quant Time: Apr 09 18:47:29 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 09 18:44:11 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	109	0.00
2		1,4-Dioxane	1.125	1.172	-4.2	112	0.01
3		Pyridine	2.523	2.764	-9.6	110	0.00
4		n-Nitrosodimethylamine	0.745	0.818	-9.8	117	0.00
5	S	2-Fluorophenol	1.906	1.983	-4.0	114	0.00
6		Aniline	2.341	2.389	-2.1	112	0.00
7	S	Phenol-d6	2.174	2.216	-1.9	112	0.00
8		2-Chlorophenol	1.368	1.409	-3.0	112	0.00
9		Benzaldehyde	1.432	1.461	-2.0	114	0.00
10	C	Phenol	2.227	2.260	-1.5	112	0.00
11		bis(2-Chloroethyl)ether	1.596	1.673	-4.8	114	0.00
12		1,3-Dichlorobenzene	1.488	1.485	0.2	111	0.00
13	C	1,4-Dichlorobenzene	1.469	1.493	-1.6	111	0.00
14		1,2-Dichlorobenzene	1.397	1.398	-0.1	110	0.00
15		Benzyl Alcohol	1.506	1.519	-0.9	110	0.00
16		2,2'-oxybis(1-Chloropropane	0.929	0.909	2.2	110	0.01
17		2-Methylphenol	1.222	1.240	-1.5	112	0.00
18		Hexachloroethane	0.637	0.642	-0.8	112	0.00
19	P	n-Nitroso-di-n-propylamine	1.338	1.390	-3.9	114	0.00
20		3+4-Methylphenols	1.600	1.613	-0.8	111	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	109	0.00
22		Acetophenone	0.705	0.731	-3.7	113	0.00
23	S	Nitrobenzene-d5	0.634	0.682	-7.6	113	0.00
24		Nitrobenzene	0.656	0.697	-6.2	114	0.00
25		Isophorone	1.058	1.104	-4.3	112	0.00
26	C	2-Nitrophenol	0.182	0.190	-4.4	107	0.00
27		2,4-Dimethylphenol	0.295	0.312	-5.8	113	0.00
28		bis(2-Chloroethoxy)methane	0.502	0.521	-3.8	111	0.00
29	C	2,4-Dichlorophenol	0.328	0.342	-4.3	112	0.00
30		1,2,4-Trichlorobenzene	0.376	0.391	-4.0	113	0.00
31		Naphthalene	1.095	1.129	-3.1	112	0.00
32		Benzoic acid	0.193	0.197	-2.1	106	0.00
33		4-Chloroaniline	0.408	0.426	-4.4	111	0.00
34	C	Hexachlorobutadiene	0.260	0.269	-3.5	113	0.00
35		Caprolactam	0.109	0.114	-4.6	104	0.00
36	C	4-Chloro-3-methylphenol	0.424	0.440	-3.8	110	0.00
37		2-Methylnaphthalene	0.743	0.761	-2.4	110	0.00
38		1-Methylnaphthalene	0.703	0.723	-2.8	112	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
40		1,2,4,5-Tetrachlorobenzene	0.674	0.717	-6.4	111	0.00
41	P	Hexachlorocyclopentadiene	0.310	0.336	-8.4	111	0.00
42	S	2,4,6-Tribromophenol	0.252	0.269	-6.7	110	0.00
43	C	2,4,6-Trichlorophenol	0.423	0.460	-8.7	109	0.00
44		2,4,5-Trichlorophenol	0.462	0.484	-4.8	106	0.00
45	S	2-Fluorobiphenyl	1.481	1.558	-5.2	110	0.00
46		1,1'-Biphenyl	1.516	1.605	-5.9	110	0.00
47		2-Chloronaphthalene	1.198	1.269	-5.9	109	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.464	0.520	-12.1	112	0.00
49		Acenaphthylene	1.853	1.964	-6.0	110	0.00
50		Dimethylphthalate	1.468	1.578	-7.5	112	0.00
51		2,6-Dinitrotoluene	0.327	0.366	-11.9	111	0.00
52	C	Acenaphthene	1.132	1.184	-4.6	109	0.00
53		3-Nitroaniline	0.298	0.332	-11.4	108	0.00
54	P	2,4-Dinitrophenol	0.179	0.192	-7.3	110	0.00
55		Dibenzofuran	1.885	1.986	-5.4	110	0.00
56	P	4-Nitrophenol	0.250	0.269	-7.6	107	0.00
57		2,4-Dinitrotoluene	0.454	0.497	-9.5	108	0.00
58		Fluorene	1.528	1.630	-6.7	111	0.00
59		2,3,4,6-Tetrachlorophenol	0.411	0.441	-7.3	113	0.00
60		Diethylphthalate	1.502	1.634	-8.8	111	0.00
61		4-Chlorophenyl-phenylether	0.802	0.840	-4.7	108	0.00
62		4-Nitroaniline	0.306	0.356	-16.3	109	0.00
63		Azobenzene	2.317	2.551	-10.1	113	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
65		4,6-Dinitro-2-methylphenol	0.134	0.138	-3.0	108	0.00
66	c	n-Nitrosodiphenylamine	0.604	0.619	-2.5	111	0.00
67		4-Bromophenyl-phenylether	0.228	0.232	-1.8	111	0.00
68		Hexachlorobenzene	0.244	0.247	-1.2	109	0.00
69		Atrazine	0.211	0.227	-7.6	112	0.00
70	C	Pentachlorophenol	0.147	0.155	-5.4	110	0.00
71		Phenanthrene	1.131	1.166	-3.1	112	0.00
72		Anthracene	1.095	1.146	-4.7	111	0.00
73		Carbazole	1.046	1.077	-3.0	108	0.00
74		Di-n-butylphthalate	1.169	1.244	-6.4	111	0.00
75	C	Fluoranthene	1.387	1.428	-3.0	110	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	108	0.00
77		Benzidine	0.366	0.387	-5.7	104	0.00
78		Pyrene	1.243	1.248	-0.4	109	0.00
79	S	Terphenyl-d14	1.018	1.027	-0.9	110	0.00
80		Butylbenzylphthalate	0.464	0.460	0.9	105	0.00
81		Benzo(a)anthracene	1.320	1.327	-0.5	108	0.00
82		3,3'-Dichlorobenzidine	0.386	0.386	0.0	103	0.00
83		Chrysene	1.284	1.270	1.1	107	0.00
84		Bis(2-ethylhexyl)phthalate	0.730	0.739	-1.2	108	0.00
85	c	Di-n-octyl phthalate	1.311	1.289	1.7	106	0.00
86	I	Perylene-d12	1.000	1.000	0.0	105	0.00
87		Indeno(1,2,3-cd)pyrene	1.529	1.573	-2.9	107	0.00
88		Benzo(b)fluoranthene	1.309	1.355	-3.5	106	0.00
89		Benzo(k)fluoranthene	1.255	1.282	-2.2	105	0.00
90	C	Benzo(a)pyrene	1.212	1.249	-3.1	106	0.00
91		Dibenzo(a,h)anthracene	1.333	1.373	-3.0	106	0.00
92		Benzo(g,h,i)perylene	1.289	1.332	-3.3	107	0.00

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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