Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080321\
Data File : BP006480.D
Acq On : 03 Aug 2021 16:25
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP080321

Quant Time: Aug 03 17:02:06 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 03 15:41:30 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	114	0.00
2		1,4-Dioxane	0.567	0.562	0.9	115	0.00
3		Pyridine	1.666	1.649	1.0	108	0.00
4		n-Nitrosodimethylamine	0.629	0.623	1.0	109	0.00
5	S	2-Fluorophenol	1.302	1.311	-0.7	112	0.00
6		Aniline	2.009	2.015	-0.3	105	0.00
7	S	Phenol-d6	1.850	1.835	0.8	106	0.00
8		2-Chlorophenol	1.409	1.406	0.2	108	0.00
9		Benzaldehyde	1.113	1.086	2.4	109	0.00
10	C	Phenol	1.855	1.840	0.8	106	0.00
11		bis(2-Chloroethyl)ether	1.408	1.396	0.9	107	0.00
12		1,3-Dichlorobenzene	1.468	1.455	0.9	110	0.00
13	C	1,4-Dichlorobenzene	1.490	1.485	0.3	111	0.00
14		1,2-Dichlorobenzene	1.430	1.412	1.3	108	0.00
15		Benzyl Alcohol	1.387	1.396	-0.6	105	0.00
16		2,2'-oxybis(1-Chloropropane	1.676	1.637	2.3	104	0.00
17		2-Methylphenol	1.171	1.177	-0.5	106	0.00
18		Hexachloroethane	0.590	0.575	2.5	109	0.00
19	P	n-Nitroso-di-n-propylamine	1.241	1.243	-0.2	103	0.00
20		3+4-Methylphenols	1.594	1.616	-1.4	105	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.530	0.529	0.2	103	0.00
23	S	Nitrobenzene-d5	0.433	0.431	0.5	103	0.00
24		Nitrobenzene	0.450	0.448	0.4	104	0.00
25		Isophorone	0.778	0.796	-2.3	103	0.00
26	C	2-Nitrophenol	0.164	0.172	-4.9	106	0.00
27		2,4-Dimethylphenol	0.274	0.281	-2.6	105	0.00
28		bis(2-Chloroethoxy)methane	0.416	0.417	-0.2	104	0.00
29	C	2,4-Dichlorophenol	0.278	0.284	-2.2	103	0.00
30		1,2,4-Trichlorobenzene	0.299	0.297	0.7	106	0.00
31		Naphthalene	1.075	1.069	0.6	104	0.00
32		Benzoic acid	0.212	0.218	-2.8	106	0.00
33		4-Chloroaniline	0.435	0.440	-1.1	102	0.00
34	C	Hexachlorobutadiene	0.174	0.175	-0.6	108	0.00
35		Caprolactam	0.100	0.108	-8.0	101	0.00
36	C	4-Chloro-3-methylphenol	0.353	0.360	-2.0	101	0.00
37		2-Methylnaphthalene	0.728	0.727	0.1	103	0.00
38		1-Methylnaphthalene	0.692	0.686	0.9	102	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
40		1,2,4,5-Tetrachlorobenzene	0.523	0.512	2.1	103	0.00
41	P	Hexachlorocyclopentadiene	0.278	0.287	-3.2	107	0.00
42	S	2,4,6-Tribromophenol	0.209	0.217	-3.8	103	0.00
43	C	2,4,6-Trichlorophenol	0.357	0.365	-2.2	103	0.00
44		2,4,5-Trichlorophenol	0.396	0.401	-1.3	102	0.00
45	S	2-Fluorobiphenyl	1.337	1.316	1.6	103	0.00
46		1,1'-Biphenyl	1.514	1.483	2.0	102	0.00
47		2-Chloronaphthalene	1.166	1.139	2.3	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080321\
Data File : BP006480.D
Acq On : 03 Aug 2021 16:25
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP080321

Quant Time: Aug 03 17:02:06 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 03 15:41:30 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)
48		2-Nitroaniline	0.411	0.422	-2.7	100	0.00
49		Acenaphthylene	1.880	1.880	0.0	102	0.00
50		Dimethylphthalate	1.456	1.440	1.1	101	0.00
51		2,6-Dinitrotoluene	0.325	0.326	-0.3	99	0.00
52	C	Acenaphthene	1.144	1.127	1.5	101	0.00
53		3-Nitroaniline	0.360	0.370	-2.8	100	0.00
54	P	2,4-Dinitrophenol	0.170	0.175	-2.9	104	0.00
55		Dibenzofuran	1.799	1.763	2.0	101	0.00
56	P	4-Nitrophenol	0.313	0.324	-3.5	98	0.00
57		2,4-Dinitrotoluene	0.436	0.450	-3.2	100	0.00
58		Fluorene	1.438	1.428	0.7	101	0.00
59		2,3,4,6-Tetrachlorophenol	0.332	0.340	-2.4	102	0.00
60		Diethylphthalate	1.529	1.540	-0.7	102	0.00
61		4-Chlorophenyl-phenylether	0.651	0.639	1.8	101	0.00
62		4-Nitroaniline	0.382	0.395	-3.4	99	0.00
63		Azobenzene	1.783	1.785	-0.1	99	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	0.122	0.123	-0.8	99	0.00
66	c	n-Nitrosodiphenylamine	0.629	0.631	-0.3	100	0.00
67		4-Bromophenyl-phenylether	0.192	0.195	-1.6	103	0.00
68		Hexachlorobenzene	0.218	0.218	0.0	102	0.00
69		Atrazine	0.196	0.205	-4.6	101	0.00
70	C	Pentachlorophenol	0.139	0.147	-5.8	101	0.00
71		Phenanthrene	1.128	1.120	0.7	100	0.00
72		Anthracene	1.090	1.101	-1.0	100	0.00
73		Carbazole	1.112	1.123	-1.0	98	0.00
74		Di-n-butylphthalate	1.267	1.331	-5.1	103	0.00
75	C	Fluoranthene	1.266	1.286	-1.6	100	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
77		Benzidine	0.500	0.588	-17.6	103	0.00
78		Pyrene	1.336	1.336	0.0	99	0.00
79	S	Terphenyl-d14	0.998	0.992	0.6	101	0.00
80		Butylbenzylphthalate	0.587	0.623	-6.1	102	0.00
81		Benzo(a)anthracene	1.307	1.292	1.1	98	0.00
82		3,3'-Dichlorobenzidine	0.343	0.355	-3.5	101	0.00
83		Chrysene	1.267	1.258	0.7	99	0.00
84		Bis(2-ethylhexyl)phthalate	0.883	0.927	-5.0	103	0.00
85	c	Di-n-octyl phthalate	1.455	1.548	-6.4	103	0.00
86	I	Perylene-d12	1.000	1.000	0.0	103	0.00
87		Indeno(1,2,3-cd)pyrene	1.450	1.457	-0.5	99	0.00
88		Benzo(b)fluoranthene	1.300	1.292	0.6	99	0.00
89		Benzo(k)fluoranthene	1.248	1.253	-0.4	99	0.00
90	C	Benzo(a)pyrene	1.166	1.181	-1.3	99	0.00
91		Dibenzo(a,h)anthracene	1.254	1.251	0.2	99	0.00
92		Benzo(g,h,i)perylene	1.202	1.196	0.5	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080321\
Data File : BP006480.D
Acq On : 03 Aug 2021 16:25
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP080321

Quant Time: Aug 03 17:02:06 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 03 15:41:30 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)

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