Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110619\
Data File : BF117703.D
Acq On : 6 Nov 2019 17:22
Operator : JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF110619

Quant Time: Nov 06 17:51:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 06 17:27:10 2019
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	100	0.00
2		1,4-Dioxane	0.551	0.538	2.4	102	0.00
3		Pyridine	1.514	1.460	3.6	101	0.00
4		n-Nitrosodimethylamine	0.667	0.649	2.7	101	0.00
5	S	2-Fluorophenol	1.105	1.073	2.9	100	0.00
6		Aniline	1.912	1.835	4.0	99	0.00
7	S	Phenol-d6	1.362	1.324	2.8	100	0.00
8		2-Chlorophenol	1.229	1.210	1.5	101	0.00
9		Benzaldehyde	0.948	0.977	-3.1	104	0.00
10	C	Phenol	1.468	1.438	2.0	102	0.00
11		bis(2-Chloroethyl)ether	1.212	1.186	2.1	101	0.00
12		1,3-Dichlorobenzene	1.425	1.406	1.3	101	0.00
13	C	1,4-Dichlorobenzene	1.426	1.400	1.8	101	0.00
14		1,2-Dichlorobenzene	1.313	1.307	0.5	101	0.00
15		Benzyl Alcohol	1.086	1.068	1.7	101	0.00
16		2,2'-oxybis(1-Chloropropane	1.917	1.861	2.9	99	0.00
17		2-Methylphenol	1.056	1.027	2.7	101	0.00
18		Hexachloroethane	0.527	0.524	0.6	103	0.00
19	P	n-Nitroso-di-n-propylamine	0.896	0.866	3.3	101	0.00
20		3+4-Methylphenols	1.257	1.245	1.0	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
22		Acetophenone	0.457	0.442	3.3	101	0.00
23	S	Nitrobenzene-d5	0.339	0.336	0.9	104	0.00
24		Nitrobenzene	0.354	0.349	1.4	102	0.00
25		Isophorone	0.648	0.611	5.7	101	0.00
26	C	2-Nitrophenol	0.164	0.164	0.0	103	0.00
27		2,4-Dimethylphenol	0.277	0.263	5.1	100	0.00
28		bis(2-Chloroethoxy)methane	0.416	0.398	4.3	101	0.00
29	C	2,4-Dichlorophenol	0.285	0.273	4.2	99	0.00
30		1,2,4-Trichlorobenzene	0.327	0.317	3.1	100	0.00
31		Naphthalene	0.914	0.894	2.2	101	0.00
32		Benzoic acid	0.219	0.217	0.9	109	0.00
33		4-Chloroaniline	0.425	0.399	6.1	98	0.00
34	C	Hexachlorobutadiene	0.187	0.183	2.1	101	0.00
35		Caprolactam	0.095	0.089	6.3	102	0.00
36	C	4-Chloro-3-methylphenol	0.288	0.274	4.9	100	0.00
37		2-Methylnaphthalene	0.622	0.606	2.6	100	0.00
38		1-Methylnaphthalene	0.591	0.575	2.7	100	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
40		1,2,4,5-Tetrachlorobenzene	0.547	0.537	1.8	100	0.00
41	P	Hexachlorocyclopentadiene	0.295	0.280	5.1	97	0.00
42	S	2,4,6-Tribromophenol	0.196	0.189	3.6	100	0.00
43	C	2,4,6-Trichlorophenol	0.398	0.382	4.0	100	0.00
44		2,4,5-Trichlorophenol	0.404	0.391	3.2	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110619\
Data File : BF117703.D
Acq On : 6 Nov 2019 17:22
Operator : JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF110619

Quant Time: Nov 06 17:51:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 06 17:27:10 2019
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)
45	S	2-Fluorobiphenyl	1.014	1.068	-5.3	101	0.00
46		1,1'-Biphenyl	1.401	1.383	1.3	101	0.00
47		2-Chloronaphthalene	1.162	1.139	2.0	100	0.00
48		2-Nitroaniline	0.364	0.354	2.7	101	0.00
49		Acenaphthylene	1.693	1.660	1.9	100	0.00
50		Dimethylphthalate	1.352	1.300	3.8	100	0.00
51		2,6-Dinitrotoluene	0.303	0.294	3.0	101	0.00
52	C	Acenaphthene	1.092	1.061	2.8	100	0.00
53		3-Nitroaniline	0.365	0.356	2.5	102	0.00
54	P	2,4-Dinitrophenol	0.104	0.104	0.0	108	0.00
55		Dibenzofuran	1.529	1.491	2.5	99	0.00
56	P	4-Nitrophenol	0.275	0.267	2.9	101	0.00
57		2,4-Dinitrotoluene	0.388	0.384	1.0	102	0.00
58		Fluorene	1.127	1.082	4.0	98	0.00
59		2,3,4,6-Tetrachlorophenol	0.336	0.325	3.3	100	0.00
60		Diethylphthalate	1.330	1.299	2.3	100	0.00
61		4-Chlorophenyl-phenylether	0.586	0.571	2.6	99	0.00
62		4-Nitroaniline	0.360	0.338	6.1	98	0.00
63		Azobenzene	1.259	1.222	2.9	100	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
65		4,6-Dinitro-2-methylphenol	0.080	0.078	2.5	102	0.00
66	c	n-Nitrosodiphenylamine	0.591	0.575	2.7	100	0.00
67		4-Bromophenyl-phenylether	0.218	0.210	3.7	99	0.00
68		Hexachlorobenzene	0.252	0.247	2.0	101	0.00
69		Atrazine	0.195	0.185	5.1	98	0.00
70	C	Pentachlorophenol	0.141	0.138	2.1	100	0.00
71		Phenanthrene	0.959	0.934	2.6	100	0.00
72		Anthracene	0.954	0.937	1.8	100	0.00
73		Carbazole	0.872	0.844	3.2	98	0.00
74		Di-n-butylphthalate	1.060	1.044	1.5	100	0.00
75	C	Fluoranthene	0.979	0.947	3.3	99	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	97	0.00
77		Benzidine	0.695	0.621	10.6	85	0.00
78		Pyrene	1.460	1.421	2.7	99	0.00
79	S	Terphenyl-d14	0.972	0.938	3.5	99	0.00
80		Butylbenzylphthalate	0.674	0.665	1.3	100	0.00
81		Benzo(a)anthracene	1.245	1.216	2.3	98	0.00
82		3,3'-Dichlorobenzidine	0.465	0.446	4.1	96	0.00
83		Chrysene	1.229	1.196	2.7	98	0.00
84		Bis(2-ethylhexyl)phthalate	0.870	0.862	0.9	100	0.00
85	c	Di-n-octyl phthalate	1.369	1.378	-0.7	101	0.00
86		Indeno(1,2,3-cd)pyrene	1.144	1.057	7.6	100	0.00
87	I	Perylene-d12	1.000	1.000	0.0	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110619\
Data File : BF117703.D
Acq On : 6 Nov 2019 17:22
Operator : JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF110619

Quant Time: Nov 06 17:51:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 06 17:27:10 2019
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)
88	Benzo(b)fluoranthene	1.260	1.207	4.2	99	0.00
89	Benzo(k)fluoranthene	1.123	1.119	0.4	96	0.00
90 C	Benzo(a)pyrene	1.092	1.039	4.9	97	0.00
91	Dibenzo(a,h)anthracene	0.979	0.916	6.4	98	0.00
92	Benzo(g,h,i)perylene	0.949	0.887	6.5	100	0.00

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