Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041819\
Data File : BF113816.D
Acq On : 18 Apr 2019 14:35
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF041819

Quant Time: Apr 18 15:00:25 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF041819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 18 13:55:51 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	99	0.00
2		1,4-Dioxane	0.599	0.597	0.3	99	0.00
3		Pyridine	1.647	1.682	-2.1	102	0.00
4		n-Nitrosodimethylamine	0.575	0.580	-0.9	99	0.00
5	S	2-Fluorophenol	1.203	1.230	-2.2	99	0.00
6		Aniline	2.129	2.148	-0.9	99	0.00
7	S	Phenol-d6	1.500	1.547	-3.1	100	0.00
8		2-Chlorophenol	1.350	1.384	-2.5	99	0.00
9		Benzaldehyde	0.892	0.897	-0.6	101	0.00
10	C	Phenol	1.681	1.738	-3.4	101	0.00
11		bis(2-Chloroethyl)ether	1.365	1.386	-1.5	99	0.00
12		1,3-Dichlorobenzene	1.500	1.534	-2.3	99	0.00
13	C	1,4-Dichlorobenzene	1.501	1.550	-3.3	101	0.00
14		1,2-Dichlorobenzene	1.383	1.443	-4.3	101	0.00
15		Benzyl Alcohol	1.215	1.265	-4.1	101	0.00
16		2,2'-oxybis(1-Chloropropane	1.656	1.685	-1.8	100	0.00
17		2-Methylphenol	1.103	1.120	-1.5	100	0.00
18		Hexachloroethane	0.538	0.548	-1.9	100	0.00
19	P	n-Nitroso-di-n-propylamine	1.001	1.007	-0.6	100	0.00
20		3+4-Methylphenols	1.325	1.391	-5.0	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.456	0.464	-1.8	99	0.00
23	S	Nitrobenzene-d5	0.328	0.347	-5.8	103	0.00
24		Nitrobenzene	0.369	0.385	-4.3	101	0.00
25		Isophorone	0.680	0.680	0.0	100	0.00
26	C	2-Nitrophenol	0.147	0.164	-11.6	110	0.00
27		2,4-Dimethylphenol	0.280	0.287	-2.5	100	0.00
28		bis(2-Chloroethoxy)methane	0.432	0.435	-0.7	99	0.00
29	C	2,4-Dichlorophenol	0.298	0.304	-2.0	99	0.00
30		1,2,4-Trichlorobenzene	0.331	0.342	-3.3	101	0.00
31		Naphthalene	0.946	0.977	-3.3	100	0.00
32		Benzoic acid	0.173	0.191	-10.4	111	0.00
33		4-Chloroaniline	0.401	0.409	-2.0	100	0.00
34	C	Hexachlorobutadiene	0.202	0.208	-3.0	100	0.00
35		Caprolactam	0.096	0.097	-1.0	99	0.00
36	C	4-Chloro-3-methylphenol	0.296	0.302	-2.0	100	0.00
37		2-Methylnaphthalene	0.628	0.642	-2.2	100	0.00
38		1-Methylnaphthalene	0.609	0.623	-2.3	99	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
40		1,2,4,5-Tetrachlorobenzene	0.635	0.652	-2.7	101	0.00
41	P	Hexachlorocyclopentadiene	0.349	0.354	-1.4	99	0.00
42	S	2,4,6-Tribromophenol	0.221	0.233	-5.4	102	0.00
43	C	2,4,6-Trichlorophenol	0.409	0.415	-1.5	101	0.00
44		2,4,5-Trichlorophenol	0.441	0.459	-4.1	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041819\
Data File : BF113816.D
Acq On : 18 Apr 2019 14:35
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF041819

Quant Time: Apr 18 15:00:25 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF041819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 18 13:55:51 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	1.311	1.263	3.7	100	0.00
46		1,1'-Biphenyl	1.513	1.566	-3.5	101	0.00
47		2-Chloronaphthalene	1.239	1.268	-2.3	100	0.00
48		2-Nitroaniline	0.322	0.355	-10.2	106	0.00
49		Acenaphthylene	1.938	1.983	-2.3	99	0.00
50		Dimethylphthalate	1.400	1.434	-2.4	100	0.00
51		2,6-Dinitrotoluene	0.309	0.329	-6.5	104	0.00
52	C	Acenaphthene	1.137	1.174	-3.3	102	0.00
53		3-Nitroaniline	0.327	0.354	-8.3	106	0.00
54	P	2,4-Dinitrophenol	0.094	0.107	-13.8	118	0.00
55		Dibenzofuran	1.703	1.766	-3.7	101	0.00
56	P	4-Nitrophenol	0.266	0.291	-9.4	105	0.00
57		2,4-Dinitrotoluene	0.383	0.421	-9.9	107	0.00
58		Fluorene	1.270	1.295	-2.0	100	0.00
59		2,3,4,6-Tetrachlorophenol	0.376	0.385	-2.4	100	0.00
60		Diethylphthalate	1.341	1.378	-2.8	101	0.00
61		4-Chlorophenyl-phenylether	0.650	0.670	-3.1	100	0.00
62		4-Nitroaniline	0.315	0.330	-4.8	104	0.00
63		Azobenzene	1.378	1.412	-2.5	100	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
65		4,6-Dinitro-2-methylphenol	0.076	0.084	-10.5	111	0.00
66	c	n-Nitrosodiphenylamine	0.606	0.612	-1.0	101	0.00
67		4-Bromophenyl-phenylether	0.235	0.238	-1.3	100	0.00
68		Hexachlorobenzene	0.258	0.262	-1.6	100	0.00
69		Atrazine	0.188	0.190	-1.1	98	0.00
70	C	Pentachlorophenol	0.146	0.154	-5.5	105	0.00
71		Phenanthrene	1.019	1.033	-1.4	100	0.00
72		Anthracene	1.026	1.050	-2.3	101	0.00
73		Carbazole	0.932	0.951	-2.0	101	0.00
74		Di-n-butylphthalate	1.037	1.054	-1.6	100	0.00
75	C	Fluoranthene	1.080	1.118	-3.5	101	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
77		Benzidine	0.579	0.659	-13.8	108	0.00
78		Pyrene	1.410	1.330	5.7	101	0.00
79	S	Terphenyl-d14	0.954	0.911	4.5	101	0.00
80		Butylbenzylphthalate	0.522	0.515	1.3	105	0.00
81		Benzo(a)anthracene	1.169	1.164	0.4	104	0.00
82		3,3'-Dichlorobenzidine	0.416	0.441	-6.0	106	0.00
83		Chrysene	1.160	1.170	-0.9	105	0.00
84		Bis(2-ethylhexyl)phthalate	0.625	0.627	-0.3	106	0.00
85	c	Di-n-octyl phthalate	0.969	1.047	-8.0	113	0.00
86		Indeno(1,2,3-cd)pyrene	0.947	0.879	7.2	109	0.00
87	I	Perylene-d12	1.000	1.000	0.0	110	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041819\
Data File : BF113816.D
Acq On : 18 Apr 2019 14:35
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF041819

Quant Time: Apr 18 15:00:25 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF041819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 18 13:55:51 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	1.247	1.287	-3.2	109	0.00
89	Benzo(k)fluoranthene	1.136	1.190	-4.8	112	0.00
90 C	Benzo(a)pyrene	1.102	1.121	-1.7	111	0.00
91	Dibenzo(a,h)anthracene	0.981	0.904	7.8	108	0.00
92	Benzo(g,h,i)perylene	0.997	0.884	11.3	108	0.00

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