Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012519\
Data File : BF112236.D
Acq On : 25 Jan 2019 16:45
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF012519

Quant Time: Jan 28 03:05:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 28 02:58:56 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	97	0.00
2		1,4-Dioxane	0.375	0.371	1.1	100	0.00
3		Pyridine	0.954	0.886	7.1	94	0.00
4		n-Nitrosodimethylamine	0.401	0.376	6.2	90	0.00
5	S	2-Fluorophenol	0.942	0.874	7.2	82	0.00
6		Aniline	1.556	1.551	0.3	93	0.00
7	S	Phenol-d6	1.308	1.297	0.8	94	0.00
8		2-Chlorophenol	1.267	1.257	0.8	96	0.00
9		Benzaldehyde	0.684	0.676	1.2	94	0.00
10	C	Phenol	1.435	1.449	-1.0	95	0.00
11		bis(2-Chloroethyl)ether	1.130	1.108	1.9	94	0.00
12		1,3-Dichlorobenzene	1.475	1.430	3.1	96	0.00
13	C	1,4-Dichlorobenzene	1.518	1.484	2.2	99	0.00
14		1,2-Dichlorobenzene	1.421	1.406	1.1	100	0.00
15		Benzyl Alcohol	1.103	1.101	0.2	99	0.00
16		2,2'-oxybis(1-Chloropropane	1.451	1.453	-0.1	99	0.00
17		2-Methylphenol	1.004	0.987	1.7	99	0.00
18		Hexachloroethane	0.533	0.540	-1.3	101	0.00
19	P	n-Nitroso-di-n-propylamine	0.902	0.888	1.6	99	0.00
20		3+4-Methylphenols	1.284	1.254	2.3	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	109	0.00
22		Acetophenone	0.487	0.477	2.1	99	0.00
23	S	Nitrobenzene-d5	0.347	0.348	-0.3	101	0.00
24		Nitrobenzene	0.347	0.345	0.6	99	0.00
25		Isophorone	0.545	0.550	-0.9	104	0.00
26	C	2-Nitrophenol	0.185	0.194	-4.9	108	0.00
27		2,4-Dimethylphenol	0.255	0.256	-0.4	107	0.00
28		bis(2-Chloroethoxy)methane	0.371	0.375	-1.1	110	0.00
29	C	2,4-Dichlorophenol	0.286	0.290	-1.4	109	0.00
30		1,2,4-Trichlorobenzene	0.328	0.330	-0.6	111	0.00
31		Naphthalene	0.935	0.927	0.9	111	0.00
32		Benzoic acid	0.146	0.135	7.5	97	0.00
33		4-Chloroaniline	0.407	0.408	-0.2	112	0.00
34	C	Hexachlorobutadiene	0.193	0.193	0.0	112	0.00
35		Caprolactam	0.101	0.100	1.0	111	0.00
36	C	4-Chloro-3-methylphenol	0.302	0.295	2.3	109	0.00
37		2-Methylnaphthalene	0.640	0.599	6.4	105	0.00
38		1-Methylnaphthalene	0.614	0.548	10.7	101	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	83	0.00
40		1,2,4,5-Tetrachlorobenzene	0.657	0.732	-11.4	98	0.00
41	P	Hexachlorocyclopentadiene	0.195	0.231	-18.5	101	0.00
42	S	2,4,6-Tribromophenol	0.233	0.264	-13.3	107	0.00
43	C	2,4,6-Trichlorophenol	0.405	0.457	-12.8	98	0.00
44		2,4,5-Trichlorophenol	0.423	0.483	-14.2	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012519\
Data File : BF112236.D
Acq On : 25 Jan 2019 16:45
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF012519

Quant Time: Jan 28 03:05:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 28 02:58:56 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	1.414	1.533	-8.4	99	0.00
46		1,1'-Biphenyl	1.748	1.916	-9.6	98	0.00
47		2-Chloronaphthalene	1.356	1.490	-9.9	98	0.00
48		2-Nitroaniline	0.370	0.403	-8.9	94	0.00
49		Acenaphthylene	1.987	2.180	-9.7	97	0.00
50		Dimethylphthalate	1.554	1.692	-8.9	98	0.00
51		2,6-Dinitrotoluene	0.348	0.384	-10.3	98	0.00
52	C	Acenaphthene	1.187	1.173	1.2	86	0.00
53		3-Nitroaniline	0.377	0.422	-11.9	99	0.00
54	P	2,4-Dinitrophenol	0.112	0.113	-0.9	88	0.00
55		Dibenzofuran	1.814	2.000	-10.3	97	0.00
56	P	4-Nitrophenol	0.199	0.229	-15.1	96	0.00
57		2,4-Dinitrotoluene	0.443	0.506	-14.2	99	0.00
58		Fluorene	1.358	1.499	-10.4	104	0.00
59		2,3,4,6-Tetrachlorophenol	0.320	0.365	-14.1	94	0.00
60		Diethylphthalate	1.542	1.720	-11.5	99	0.00
61		4-Chlorophenyl-phenylether	0.738	0.810	-9.8	98	0.00
62		4-Nitroaniline	0.370	0.382	-3.2	92	0.00
63		Azobenzene	1.356	1.542	-13.7	97	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	81	0.00
65		4,6-Dinitro-2-methylphenol	0.112	0.128	-14.3	95	0.00
66	c	n-Nitrosodiphenylamine	0.674	0.745	-10.5	95	0.00
67		4-Bromophenyl-phenylether	0.235	0.250	-6.4	90	0.00
68		Hexachlorobenzene	0.252	0.286	-13.5	98	0.00
69		Atrazine	0.217	0.255	-17.5	99	0.00
70	C	Pentachlorophenol	0.098	0.116	-18.4	103	0.00
71		Phenanthrene	1.040	1.059	-1.8	86	0.00
72		Anthracene	1.036	1.108	-6.9	100	0.00
73		Carbazole	1.022	1.071	-4.8	89	0.00
74		Di-n-butylphthalate	1.268	1.408	-11.0	95	0.00
75	C	Fluoranthene	1.115	1.307	-17.2	93	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	84	0.00
77		Benzidine	0.550	0.498	9.5	74	0.00
78		Pyrene	1.385	1.423	-2.7	94	0.00
79	S	Terphenyl-d14	1.062	1.058	0.4	95	0.00
80		Butylbenzylphthalate	0.670	0.683	-1.9	93	0.00
81		Benzo(a)anthracene	1.176	1.156	1.7	87	0.00
82		3,3'-Dichlorobenzidine	0.440	0.424	3.6	86	0.00
83		Chrysene	1.122	1.113	0.8	91	0.00
84		Bis(2-ethylhexyl)phthalate	0.951	0.947	0.4	90	0.00
85	c	Di-n-octyl phthalate	1.463	1.429	2.3	88	0.00
86		Indeno(1,2,3-cd)pyrene	0.889	0.874	1.7	103	0.00
87	I	Perylene-d12	1.000	1.000	0.0	82	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012519\
Data File : BF112236.D
Acq On : 25 Jan 2019 16:45
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF012519

Quant Time: Jan 28 03:05:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 28 02:58:56 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	1.196	1.254	-4.8	88	0.00
89	Benzo(k)fluoranthene	1.146	1.125	1.8	85	0.00
90 C	Benzo(a)pyrene	1.076	1.077	-0.1	86	0.00
91	Dibenzo(a,h)anthracene	0.867	0.947	-9.2	104	0.00
92	Benzo(g,h,i)perylene	0.818	0.899	-9.9	114	0.00

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