Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF061818\
Data File : BF106525.D
Acq On : 18 Jun 2018 9:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 19 09:38:16 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 15 11:11:47 2018
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	59	0.00
2		1,4-Dioxane	0.612	0.569	7.0	57	0.03
3		Pyridine	1.704	1.579	7.3	57	0.02
4		n-Nitrosodimethylamine	0.781	0.697	10.8	53	0.02
5	S	2-Fluorophenol	1.182	1.143	3.3	59	0.00
6		Aniline	2.180	1.927	11.6	54	0.00
7	S	Phenol-d6	1.493	1.399	6.3	58	0.00
8		2-Chlorophenol	1.271	1.249	1.7	60	0.00
9		Benzaldehyde	1.132	1.022	9.7	56	0.00
10	C	Phenol	1.630	1.622	0.5	62	0.00
11		bis(2-Chloroethyl)ether	1.395	1.352	3.1	60	0.00
12		1,3-Dichlorobenzene	1.496	1.467	1.9	60	0.00
13	C	1,4-Dichlorobenzene	1.518	1.476	2.8	60	0.00
14		1,2-Dichlorobenzene	1.429	1.370	4.1	59	0.00
15		Benzyl Alcohol	1.279	1.207	5.6	56	0.00
16		2,2'-oxybis(1-Chloropropane	2.000	1.790	10.5	54	0.00
17		2-Methylphenol	1.134	1.062	6.3	57	0.00
18		Hexachloroethane	0.541	0.533	1.5	59	0.00
19	P	n-Nitroso-di-n-propylamine	1.120	0.938	16.3	52	0.00
20		3+4-Methylphenols	1.450	1.250	13.8	53	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	59	0.00
22		Acetophenone	0.504	0.436	13.5	54	0.00
23	S	Nitrobenzene-d5	0.340	0.345	-1.5	60	0.00
24		Nitrobenzene	0.370	0.357	3.5	57	0.00
25		Isophorone	0.664	0.616	7.2	57	0.00
26	C	2-Nitrophenol	0.143	0.170	-18.9	69	0.00
27		2,4-Dimethylphenol	0.281	0.264	6.0	56	0.00
28		bis(2-Chloroethoxy)methane	0.431	0.416	3.5	59	0.00
29	C	2,4-Dichlorophenol	0.271	0.276	-1.8	60	0.00
30		1,2,4-Trichlorobenzene	0.304	0.297	2.3	59	0.00
31		Naphthalene	0.904	0.892	1.3	61	0.00
32		Benzoic acid	0.187	0.153	18.2	49#	0.02
33		4-Chloroaniline	0.401	0.385	4.0	58	0.00
34	C	Hexachlorobutadiene	0.195	0.197	-1.0	61	0.00
35		Caprolactam	0.092	0.090	2.2	59	0.02
36	C	4-Chloro-3-methylphenol	0.300	0.291	3.0	59	0.00
37		2-Methylnaphthalene	0.616	0.599	2.8	60	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	61	0.00
39		1,2,4,5-Tetrachlorobenzene	0.649	0.633	2.5	60	0.00
40	P	Hexachlorocyclopentadiene	0.358	0.351	2.0	56	0.00
41	S	2,4,6-Tribromophenol	0.192	0.224	-16.7	67	0.00
42	C	2,4,6-Trichlorophenol	0.414	0.427	-3.1	62	0.00
43		2,4,5-Trichlorophenol	0.446	0.453	-1.6	60	0.00
44	S	2-Fluorobiphenyl	1.173	1.167	0.5	62	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF061818\
Data File : BF106525.D
Acq On : 18 Jun 2018 9:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 19 09:38:16 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 15 11:11:47 2018
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		1,1'-Biphenyl	1.549	1.492	3.7	61	0.00
46		2-Chloronaphthalene	1.240	1.186	4.4	61	0.00
47		2-Nitroaniline	0.390	0.415	-6.4	60	0.00
48		Acenaphthylene	1.898	1.865	1.7	62	0.00
49		Dimethylphthalate	1.429	1.448	-1.3	64	0.01
50		2,6-Dinitrotoluene	0.292	0.306	-4.8	61	0.00
51	C	Acenaphthene	1.225	1.131	7.7	59	0.00
52		3-Nitroaniline	0.346	0.368	-6.4	63	0.00
53	P	2,4-Dinitrophenol	0.108	0.132	-22.2	75	0.00
54		Dibenzofuran	1.650	1.622	1.7	62	0.00
55	P	4-Nitrophenol	0.266	0.249	6.4	55	0.00
56		2,4-Dinitrotoluene	0.367	0.417	-13.6	65	0.01
57		Fluorene	1.307	1.278	2.2	63	0.00
58		2,3,4,6-Tetrachlorophenol	0.354	0.355	-0.3	60	0.00
59		Diethylphthalate	1.418	1.386	2.3	61	0.00
60		4-Chlorophenyl-phenylether	0.688	0.685	0.4	63	0.00
61		4-Nitroaniline	0.337	0.360	-6.8	64	0.00
62		Azobenzene	1.479	1.382	6.6	60	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	62	0.00
64		4,6-Dinitro-2-methylphenol	0.098	0.120	-22.4	75	0.00
65	c	n-Nitrosodiphenylamine	0.672	0.626	6.8	60	0.00
66		4-Bromophenyl-phenylether	0.227	0.230	-1.3	64	0.00
67		Hexachlorobenzene	0.235	0.241	-2.6	65	0.00
68		Atrazine	0.207	0.213	-2.9	66	0.00
69	C	Pentachlorophenol	0.144	0.125	13.2	51	0.00
70		Phenanthrene	0.979	0.934	4.6	64	0.00
71		Anthracene	0.981	0.948	3.4	64	0.00
72		Carbazole	0.906	0.872	3.8	63	0.00
73		Di-n-butylphthalate	1.008	1.070	-6.2	68	0.00
74	C	Fluoranthene	1.045	1.019	2.5	64	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	62	0.00
76		Benzidine	0.666	0.629	5.6	59	0.00
77		Pyrene	1.252	1.269	-1.4	65	0.00
78	S	Terphenyl-d14	0.805	0.798	0.9	66	0.00
79		Butylbenzylphthalate	0.469	0.547	-16.6	65	0.00
80		Benzo(a)anthracene	1.190	1.157	2.8	62	0.00
81		3,3'-Dichlorobenzidine	0.402	0.410	-2.0	61	0.00
82		Chrysene	1.087	1.042	4.1	61	0.01
83		Bis(2-ethylhexyl)phthalate	0.672	0.693	-3.1	60	0.00
84	c	Di-n-octyl phthalate	0.969	1.093	-12.8	65	0.02
85		Indeno(1,2,3-cd)pyrene	0.867	0.831	4.2	60	0.02
86	I	Perylene-d12	1.000	1.000	0.0	61	0.02
87		Benzo(b)fluoranthene	1.209	1.194	1.2	61	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF061818\
Data File : BF106525.D
Acq On : 18 Jun 2018 9:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 19 09:38:16 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 15 11:11:47 2018
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(k)fluoranthene	1.197	1.192	0.4	60	0.02
89 C	Benzo(a)pyrene	1.087	1.077	0.9	60	0.02
90	Dibenzo(a,h)anthracene	0.907	0.906	0.1	59	0.03
91	Benzo(g,h,i)perylene	0.881	0.882	-0.1	60	0.02

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