Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091719\
Data File : BF117109.D
Acq On : 17 Sep 2019 16:24
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 18 12:00:16 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 13 16:23:17 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	101	0.00
2		1,4-Dioxane	0.536	0.414	22.8	83	0.00
3		Pyridine	1.467	1.047	28.6#	74	0.02
4		n-Nitrosodimethylamine	0.504	0.346	31.3#	74	0.00
5	S	2-Fluorophenol	1.281	1.130	11.8	92	0.00
6		Aniline	2.132	1.999	6.2	99	0.00
7	S	Phenol-d6	1.656	1.565	5.5	100	0.01
8		2-Chlorophenol	1.382	1.313	5.0	100	0.00
9		Benzaldehyde	0.904	0.875	3.2	103	0.00
10	C	Phenol	1.825	1.703	6.7	98	0.00
11		bis(2-Chloroethyl)ether	1.341	1.280	4.5	103	0.00
12		1,3-Dichlorobenzene	1.551	1.482	4.4	101	0.00
13	C	1,4-Dichlorobenzene	1.583	1.514	4.4	100	0.00
14		1,2-Dichlorobenzene	1.473	1.402	4.8	100	0.00
15		Benzyl Alcohol	1.270	1.235	2.8	101	0.00
16		2,2'-oxybis(1-Chloropropane	1.325	1.249	5.7	100	0.00
17		2-Methylphenol	1.158	1.089	6.0	101	0.00
18		Hexachloroethane	0.609	0.580	4.8	101	0.00
19	P	n-Nitroso-di-n-propylamine	1.008	0.996	1.2	106	0.00
20		3+4-Methylphenols	1.443	1.385	4.0	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
22		Acetophenone	0.508	0.479	5.7	104	0.00
23	S	Nitrobenzene-d5	0.381	0.363	4.7	102	0.00
24		Nitrobenzene	0.376	0.353	6.1	102	0.00
25		Isophorone	0.674	0.647	4.0	107	0.00
26	C	2-Nitrophenol	0.161	0.166	-3.1	110	0.00
27		2,4-Dimethylphenol	0.256	0.242	5.5	103	0.00
28		bis(2-Chloroethoxy)methane	0.415	0.396	4.6	106	0.00
29	C	2,4-Dichlorophenol	0.268	0.258	3.7	102	0.00
30		1,2,4-Trichlorobenzene	0.290	0.274	5.5	103	0.00
31		Naphthalene	0.962	0.910	5.4	102	0.00
32		Benzoic acid	0.180	0.174	3.3	108	0.01
33		4-Chloroaniline	0.427	0.401	6.1	101	0.00
34	C	Hexachlorobutadiene	0.164	0.162	1.2	106	0.00
35		Caprolactam	0.091	0.088	3.3	106	0.01
36	C	4-Chloro-3-methylphenol	0.313	0.302	3.5	106	0.00
37		2-Methylnaphthalene	0.648	0.630	2.8	106	0.00
38		1-Methylnaphthalene	0.607	0.589	3.0	106	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
40		1,2,4,5-Tetrachlorobenzene	0.516	0.488	5.4	106	0.00
41	P	Hexachlorocyclopentadiene	0.194	0.234	-20.6	141	0.00
42	S	2,4,6-Tribromophenol	0.167	0.166	0.6	110	0.00
43	C	2,4,6-Trichlorophenol	0.350	0.332	5.1	106	0.00
44		2,4,5-Trichlorophenol	0.374	0.343	8.3	104	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091719\
Data File : BF117109.D
Acq On : 17 Sep 2019 16:24
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 18 12:00:16 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 13 16:23:17 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	1.279	1.224	4.3	107	0.00
46		1,1'-Biphenyl	1.569	1.496	4.7	107	0.00
47		2-Chloronaphthalene	1.238	1.165	5.9	106	0.00
48		2-Nitroaniline	0.397	0.388	2.3	111	0.00
49		Acenaphthylene	1.855	1.769	4.6	106	0.00
50		Dimethylphthalate	1.416	1.369	3.3	108	0.00
51		2,6-Dinitrotoluene	0.288	0.289	-0.3	112	0.00
52	C	Acenaphthene	1.099	1.026	6.6	104	0.00
53		3-Nitroaniline	0.364	0.348	4.4	106	0.00
54	P	2,4-Dinitrophenol	0.102	0.113	-10.8	130	0.00
55		Dibenzofuran	1.622	1.543	4.9	106	0.00
56	P	4-Nitrophenol	0.236	0.250	-5.9	116	0.01
57		2,4-Dinitrotoluene	0.374	0.390	-4.3	114	0.00
58		Fluorene	1.307	1.264	3.3	108	0.00
59		2,3,4,6-Tetrachlorophenol	0.287	0.268	6.6	102	0.00
60		Diethylphthalate	1.393	1.369	1.7	109	0.00
61		4-Chlorophenyl-phenylether	0.565	0.568	-0.5	110	0.00
62		4-Nitroaniline	0.378	0.365	3.4	106	0.00
63		Azobenzene	1.423	1.348	5.3	106	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
65		4,6-Dinitro-2-methylphenol	0.084	0.090	-7.1	121	0.00
66	c	n-Nitrosodiphenylamine	0.691	0.663	4.1	108	0.00
67		4-Bromophenyl-phenylether	0.204	0.198	2.9	108	0.00
68		Hexachlorobenzene	0.223	0.213	4.5	108	0.00
69		Atrazine	0.170	0.173	-1.8	104	0.00
70	C	Pentachlorophenol	0.105	0.094	10.5	99	0.00
71		Phenanthrene	1.136	1.075	5.4	105	0.00
72		Anthracene	1.160	1.110	4.3	105	0.00
73		Carbazole	1.101	1.033	6.2	102	0.00
74		Di-n-butylphthalate	1.310	1.320	-0.8	110	0.00
75	C	Fluoranthene	1.167	1.095	6.2	102	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	94	0.00
77		Benzidine	0.644	0.575	10.7	88	0.00
78		Pyrene	1.678	1.607	4.2	101	0.00
79	S	Terphenyl-d14	1.070	1.062	0.7	104	0.00
80		Butylbenzylphthalate	0.793	0.808	-1.9	106	0.00
81		Benzo(a)anthracene	1.336	1.276	4.5	97	0.00
82		3,3'-Dichlorobenzidine	0.431	0.408	5.3	95	0.00
83		Chrysene	1.303	1.229	5.7	96	0.00
84		Bis(2-ethylhexyl)phthalate	1.022	1.088	-6.5	111	0.00
85	c	Di-n-octyl phthalate	1.609	1.723	-7.1	108	0.00
86		Indeno(1,2,3-cd)pyrene	0.951	0.786	17.4	93	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\BF091719\
Data File : BF117109.D
Acq On : 17 Sep 2019 16:24
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 18 12:00:16 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 13 16:23:17 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	1.211	1.225	-1.2	92	0.00
89	Benzo(k)fluoranthene	1.148	1.160	-1.0	85	0.00
90 C	Benzo(a)pyrene	1.063	1.048	1.4	87	0.00
91	Dibenzo(a,h)anthracene	0.847	0.829	2.1	95	0.00
92	Benzo(g,h,i)perylene	0.844	0.783	7.2	92	0.01

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