



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:45 AM UTC

PDB ID : 1AXM / pdb_00001axm
Title : HEPARIN-LINKED BIOLOGICALLY-ACTIVE DIMER OF FIBROBLAST GROWTH FACTOR
Authors : DiGabriele, A.D.; Lax, I.; Chen, D.I.; Svahn, C.M.; Jaye, M.; Schlessinger, J.; Hendrickson, W.A.
Deposited on : 1997-10-16
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

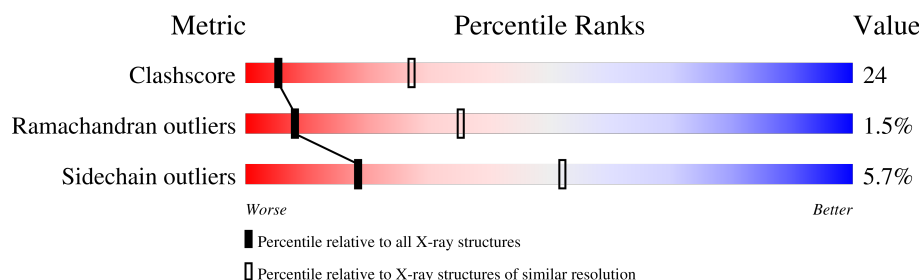
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	135	61% 34% . .
1	B	135	48% 40% . . 6%
1	C	135	47% 40% 7% . 5%
1	D	135	48% 39% 7% . 6%
1	E	135	52% 39% 5% .
1	F	135	53% 41% . .
2	G	5	100%
2	H	5	100%

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Mol	Chain	Length	Quality of chain
3	I	5	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	IDS	I	5	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6463 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

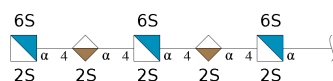
- Molecule 1 is a protein called ACIDIC FIBROBLAST GROWTH FACTOR.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	130	Total	C	N	O	S	Se	0	0	0
			1029	649	180	196	3	1			
1	B	127	Total	C	N	O	S	Se	0	0	0
			1011	638	177	192	3	1			
1	C	128	Total	C	N	O	S	Se	0	0	0
			1018	643	178	193	3	1			
1	D	127	Total	C	N	O	S	Se	0	0	0
			1011	638	177	192	3	1			
1	E	129	Total	C	N	O	S	Se	0	0	0
			1028	649	180	195	3	1			
1	F	129	Total	C	N	O	S	Se	0	0	0
			1028	649	180	195	3	1			

There are 6 discrepancies between the modelled and reference sequences:

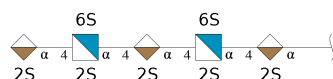
Chain	Residue	Modelled	Actual	Comment	Reference
A	67	MSE	MET	modified residue	UNP P05230
B	67	MSE	MET	modified residue	UNP P05230
C	67	MSE	MET	modified residue	UNP P05230
D	67	MSE	MET	modified residue	UNP P05230
E	67	MSE	MET	modified residue	UNP P05230
F	67	MSE	MET	modified residue	UNP P05230

- Molecule 2 is an oligosaccharide called 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	5	Total	C	N	O	S	0	0	0
			90	30	3	49	8			
2	H	5	Total	C	N	O	S	0	0	0
			90	30	3	49	8			

- Molecule 3 is an oligosaccharide called 2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	5	Total	C	N	O	S	0	0	0
			87	30	2	48	7			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	25	Total	O	0	0
			25	25		
4	B	10	Total	O	0	0
			10	10		
4	C	7	Total	O	0	0
			7	7		
4	D	13	Total	O	0	0
			13	13		
4	E	12	Total	O	0	0
			12	12		
4	F	4	Total	O	0	0
			4	4		

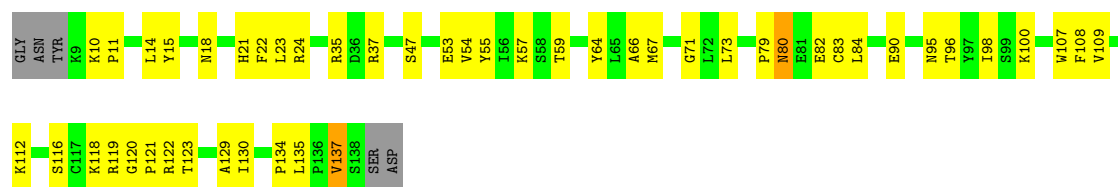
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

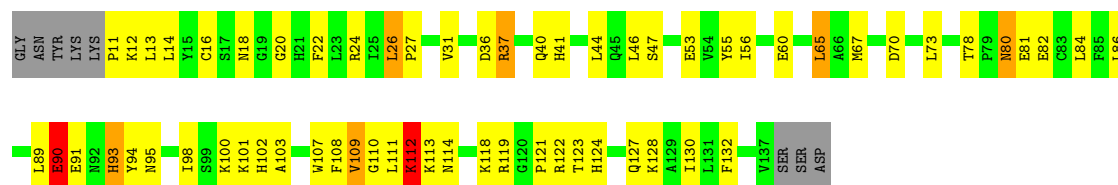
• Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain A: 



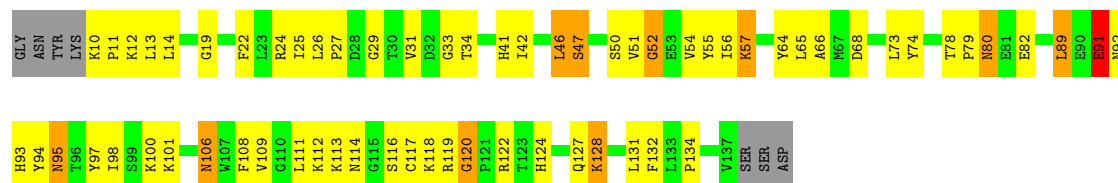
• Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain B: 



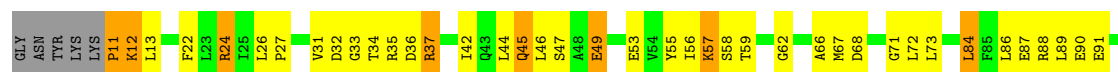
• Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

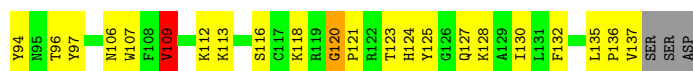
Chain C: 



• Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

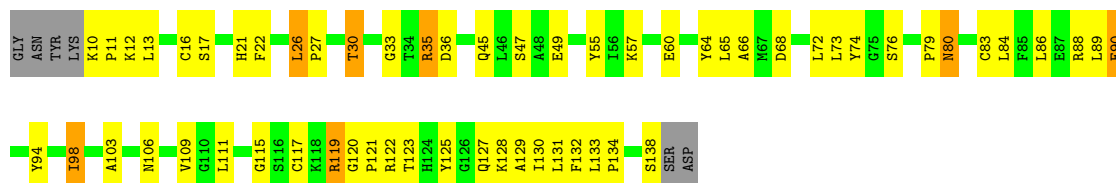
Chain D: 





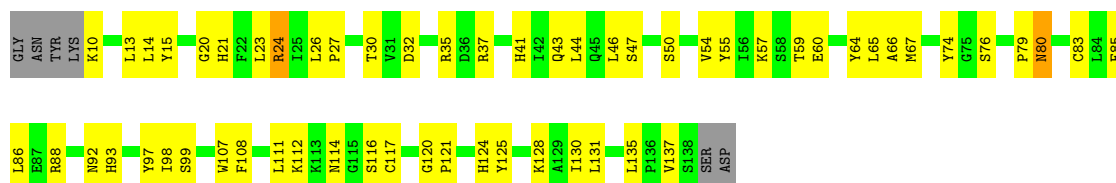
- Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain E: 52% 39% 5%



- Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain F: 53% 41%



- Molecule 2: 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose

Chain G: 100%



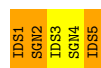
- Molecule 2: 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose

Chain H: 100%



- Molecule 3: 2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid

Chain I: 40% 60%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	87.69Å 158.05Å 190.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.00 – 3.00	Depositor
% Data completeness (in resolution range)	88.6 (14.00-3.00)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.217 , 0.304	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6463	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: IDS, SGN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.92	0/1051	1.01	2/1418 (0.1%)
1	B	0.80	0/1032	1.08	8/1391 (0.6%)
1	C	0.80	0/1040	1.01	3/1403 (0.2%)
1	D	0.80	0/1032	1.06	9/1391 (0.6%)
1	E	0.83	0/1050	1.07	4/1415 (0.3%)
1	F	0.76	0/1050	1.08	6/1415 (0.4%)
All	All	0.82	0/6255	1.05	32/8433 (0.4%)

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	12	LYS	N-CA-C	7.67	121.75	110.17
1	F	10	LYS	CA-C-N	7.36	127.41	119.90
1	F	10	LYS	C-N-CA	7.36	127.41	119.90
1	D	11	PRO	N-CA-CB	7.07	110.78	103.00
1	B	11	PRO	N-CA-CB	7.06	110.77	103.00
1	F	60	GLU	N-CA-C	6.64	118.17	111.07
1	D	24	ARG	N-CA-C	6.56	119.09	108.34
1	B	112	LYS	N-CA-C	-6.21	102.33	110.53
1	B	65	LEU	N-CA-C	-6.17	101.35	110.48
1	F	59	THR	N-CA-C	-6.13	104.52	111.14
1	C	91	GLU	N-CA-C	-6.11	102.31	110.55
1	F	93	HIS	N-CA-C	5.96	119.91	111.92
1	F	24	ARG	N-CA-C	5.95	118.14	108.26
1	D	120	GLY	CA-C-N	5.82	125.50	119.56
1	D	120	GLY	C-N-CA	5.82	125.50	119.56
1	A	109	VAL	N-CA-C	-5.68	100.40	108.58
1	B	90	GLU	N-CA-C	-5.62	102.97	110.55
1	D	37	ARG	N-CA-C	-5.57	105.31	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	26	LEU	CA-C-N	5.54	125.21	119.56
1	E	26	LEU	C-N-CA	5.54	125.21	119.56
1	A	90	GLU	N-CA-C	-5.39	103.28	110.55
1	B	26	LEU	CA-C-N	5.38	125.02	119.05
1	B	26	LEU	C-N-CA	5.38	125.02	119.05
1	D	128	LYS	N-CA-C	-5.37	105.59	111.82
1	E	35	ARG	N-CA-C	-5.32	105.53	112.23
1	D	136	PRO	N-CA-C	5.28	118.61	111.22
1	B	93	HIS	N-CA-C	5.26	117.07	110.91
1	E	128	LYS	N-CA-C	-5.23	105.75	111.82
1	B	109	VAL	N-CA-C	-5.17	102.18	109.21
1	D	109	VAL	N-CA-C	-5.10	101.23	108.58
1	C	78	THR	CA-C-N	5.04	126.15	119.84
1	C	78	THR	C-N-CA	5.04	126.15	119.84

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1029	0	1003	34	0
1	B	1011	0	989	50	0
1	C	1018	0	996	57	0
1	D	1011	0	989	62	0
1	E	1028	0	1012	47	0
1	F	1028	0	1012	42	0
2	G	90	0	31	6	0
2	H	90	0	31	10	0
3	I	87	0	30	6	0
4	A	25	0	0	2	0
4	B	10	0	0	0	0
4	C	7	0	0	2	0
4	D	13	0	0	0	0
4	E	12	0	0	1	0
4	F	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6463	0	6093	300	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 24.

All (300) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:ASN:HB2	1:B:128:LYS:HG2	1.40	1.03
1:C:47:SER:HB2	1:C:57:LYS:HE2	1.54	0.90
1:D:118:LYS:HG2	1:D:123:THR:HG22	1.53	0.89
1:A:64:TYR:CE1	1:A:79:PRO:HG3	2.13	0.82
1:E:86:LEU:HB3	1:E:88:ARG:HH21	1.46	0.81
1:D:22:PHE:HE1	1:D:135:LEU:HD23	1.47	0.79
1:E:80:ASN:ND2	1:E:83:CYS:SG	2.55	0.79
1:C:73:LEU:HD21	1:C:109:VAL:HG13	1.65	0.77
1:E:57:LYS:HB3	1:E:64:TYR:CD1	2.19	0.77
1:C:24:ARG:HH11	1:C:26:LEU:HD11	1.49	0.76
1:C:46:LEU:HD13	1:C:54:VAL:HG11	1.67	0.76
1:B:37:ARG:HG2	1:B:37:ARG:HH11	1.50	0.75
1:C:24:ARG:NH1	1:C:26:LEU:HD11	2.02	0.74
1:B:73:LEU:HD21	1:B:109:VAL:HG13	1.69	0.73
1:F:35:ARG:HG2	1:F:35:ARG:HH11	1.54	0.73
1:B:14:LEU:HB3	1:B:132:PHE:HD2	1.54	0.71
1:D:123:THR:HA	1:D:127:GLN:OE1	1.90	0.71
1:D:35:ARG:HG3	1:D:35:ARG:HH11	1.53	0.71
1:D:113:LYS:NZ	3:I:5:IDS:H2	2.06	0.70
1:D:90:GLU:HG2	1:D:96:THR:OG1	1.92	0.70
1:C:10:LYS:N	1:C:11:PRO:HD2	2.06	0.70
1:E:10:LYS:N	1:E:138:SER:HG	1.88	0.69
1:D:112:LYS:HD3	1:D:116:SER:HB2	1.75	0.69
1:E:111:LEU:HD23	1:E:117:CYS:HA	1.73	0.69
1:D:106:ASN:ND2	1:D:124:HIS:HA	2.08	0.69
1:F:43:GLN:HE22	1:F:137:VAL:HG21	1.60	0.67
1:C:12:LYS:HD3	1:C:134:PRO:HB2	1.76	0.67
1:A:96:THR:HG22	1:A:130:ILE:O	1.95	0.67
1:D:13:LEU:HD21	1:D:42:ILE:HG13	1.77	0.66
1:A:118:LYS:HD3	1:A:123:THR:HG22	1.76	0.66
1:D:97:TYR:O	1:D:109:VAL:HG23	1.96	0.66
1:C:106:ASN:ND2	1:C:108:PHE:HE1	1.93	0.66
1:E:74:TYR:HE1	1:E:76:SER:HB2	1.59	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:ALA:HB2	1:A:83:CYS:SG	2.36	0.66
1:C:118:LYS:HD2	2:H:4:IDS:O1S	1.96	0.65
1:D:22:PHE:CE1	1:D:135:LEU:HD23	2.31	0.65
1:F:35:ARG:HG2	1:F:35:ARG:NH1	2.11	0.65
1:F:66:ALA:HB3	1:F:74:TYR:CE1	2.32	0.64
1:D:35:ARG:HG3	1:D:35:ARG:NH1	2.10	0.64
1:F:15:TYR:HB2	1:F:135:LEU:HD22	1.79	0.64
1:D:24:ARG:HD2	1:D:32:ASP:OD1	1.97	0.64
1:F:65:LEU:HG	1:F:85:PHE:HE2	1.63	0.64
1:C:106:ASN:ND2	1:C:108:PHE:CE1	2.66	0.64
1:E:94:TYR:CD2	1:E:131:LEU:HB3	2.33	0.64
1:F:74:TYR:HE1	1:F:76:SER:HB2	1.62	0.64
1:A:21:HIS:CD2	1:A:35:ARG:HG3	2.33	0.63
1:B:14:LEU:HB3	1:B:132:PHE:CD2	2.33	0.63
1:B:122:ARG:HH22	2:G:1:SGN:H4	1.63	0.63
1:C:66:ALA:HB3	1:C:74:TYR:CE1	2.33	0.63
1:F:14:LEU:HG	1:F:44:LEU:HD12	1.81	0.63
1:E:12:LYS:HE2	1:E:134:PRO:HB2	1.81	0.63
1:E:57:LYS:HB3	1:E:64:TYR:CE1	2.35	0.62
1:B:37:ARG:HH11	1:B:37:ARG:CG	2.11	0.62
1:B:16:CYS:O	1:B:20:GLY:HA2	2.00	0.62
1:B:37:ARG:O	1:B:37:ARG:HD3	2.01	0.61
1:D:12:LYS:NZ	1:D:46:LEU:HD22	2.16	0.61
1:D:113:LYS:HZ2	3:I:5:IDS:H2	1.64	0.60
1:E:90:GLU:HA	4:E:654:HOH:O	2.01	0.60
1:F:57:LYS:HB3	1:F:64:TYR:CD1	2.37	0.59
1:B:114:ASN:HB3	2:G:3:SGN:S2	2.42	0.59
1:C:91:GLU:HB2	1:C:94:TYR:HD2	1.67	0.59
1:D:86:LEU:HD22	1:D:88:ARG:NH2	2.17	0.59
1:A:80:ASN:HD22	1:A:82:GLU:N	2.01	0.59
1:B:110:GLY:O	1:B:111:LEU:HD23	2.02	0.59
1:D:45:GLN:NE2	1:D:59:THR:HG22	2.18	0.59
1:B:80:ASN:HD21	1:B:82:GLU:HB2	1.67	0.58
1:E:13:LEU:O	1:E:134:PRO:HA	2.04	0.58
1:E:47:SER:HB2	1:E:55:TYR:HB2	1.86	0.58
1:C:106:ASN:HD21	1:C:108:PHE:HE1	1.50	0.58
1:C:128:LYS:HA	1:C:131:LEU:HD12	1.85	0.58
1:E:66:ALA:HB3	1:E:74:TYR:CE1	2.39	0.58
1:E:98:ILE:HD11	1:E:103:ALA:HA	1.86	0.57
2:H:4:IDS:O6A	2:H:5:SGN:H5	2.05	0.57
1:B:40:GLN:O	1:B:60:GLU:HG3	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:26:LEU:HD12	1:F:30:THR:HB	1.87	0.57
1:F:120:GLY:N	1:F:121:PRO:HD2	2.19	0.57
1:F:65:LEU:HG	1:F:85:PHE:CE2	2.40	0.57
1:A:95:ASN:HD22	1:A:134:PRO:HG2	1.70	0.57
1:B:114:ASN:HB3	2:G:3:SGN:O6S	2.05	0.57
1:B:123:THR:HA	1:B:127:GLN:OE1	2.05	0.56
1:C:14:LEU:HB3	1:C:132:PHE:CD2	2.40	0.56
1:D:57:LYS:HG2	1:D:58:SER:O	2.05	0.56
1:D:120:GLY:N	1:D:121:PRO:HD2	2.21	0.56
1:C:27:PRO:HD3	1:C:41:HIS:CE1	2.41	0.56
1:F:85:PHE:HD1	1:F:99:SER:HA	1.70	0.56
1:C:55:TYR:CE2	1:C:79:PRO:HB3	2.42	0.55
1:E:80:ASN:ND2	1:E:80:ASN:H	2.04	0.55
1:B:26:LEU:HB3	1:B:27:PRO:HD2	1.89	0.55
1:A:53:GLU:OE1	1:A:84:LEU:HD21	2.07	0.55
1:A:57:LYS:HB3	4:A:620:HOH:O	2.06	0.55
1:E:123:THR:HA	1:E:127:GLN:OE1	2.07	0.55
1:B:80:ASN:ND2	1:B:82:GLU:HB2	2.21	0.55
1:C:10:LYS:N	1:C:11:PRO:CD	2.70	0.55
1:F:24:ARG:NE	1:F:26:LEU:HD21	2.22	0.55
1:D:67:MSE:HE2	1:D:71:GLY:O	2.07	0.54
1:B:47:SER:HB2	1:B:55:TYR:HB2	1.88	0.54
1:B:119:ARG:HG2	1:B:121:PRO:HD2	1.90	0.54
1:C:57:LYS:HD2	1:C:64:TYR:CE2	2.43	0.54
1:D:106:ASN:ND2	1:D:124:HIS:CD2	2.76	0.54
1:E:64:TYR:CE2	1:E:79:PRO:HB3	2.43	0.54
1:A:37:ARG:HH22	1:A:137:VAL:HG23	1.73	0.54
1:C:14:LEU:HB3	1:C:132:PHE:HD2	1.72	0.54
1:F:37:ARG:HH11	1:F:37:ARG:HG2	1.72	0.54
1:A:80:ASN:HD22	1:A:82:GLU:H	1.56	0.53
1:D:107:TRP:NE1	1:D:121:PRO:HA	2.23	0.53
1:D:106:ASN:HD21	1:D:125:TYR:H	1.54	0.53
1:B:53:GLU:HB3	1:B:84:LEU:HD11	1.91	0.53
1:D:49:GLU:HB2	1:D:55:TYR:CE1	2.43	0.52
1:D:24:ARG:NH2	1:D:34:THR:HB	2.24	0.52
1:C:50:SER:C	1:C:52:GLY:H	2.17	0.52
1:C:47:SER:CB	1:C:57:LYS:HE2	2.35	0.52
1:C:80:ASN:HD22	1:C:80:ASN:C	2.17	0.52
1:E:21:HIS:ND1	1:E:115:GLY:HA2	2.24	0.52
1:D:57:LYS:HG3	1:D:62:GLY:O	2.10	0.52
1:B:89:LEU:HA	1:B:94:TYR:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:49:GLU:HB2	1:D:55:TYR:HE1	1.75	0.52
1:D:106:ASN:HD21	1:D:124:HIS:HA	1.75	0.52
1:D:113:LYS:HD2	3:I:5:IDS:O3	2.10	0.52
2:G:1:SGN:H61	2:G:2:IDS:H2	1.91	0.52
1:E:106:ASN:HD21	1:E:125:TYR:H	1.57	0.52
1:B:24:ARG:HG3	1:B:41:HIS:HB2	1.91	0.52
1:C:95:ASN:HB3	1:C:97:TYR:CE1	2.45	0.52
1:F:24:ARG:HB3	1:F:32:ASP:OD1	2.10	0.52
1:F:64:TYR:CE2	1:F:79:PRO:HD3	2.45	0.51
1:A:67:MSE:HE2	1:A:71:GLY:O	2.11	0.51
1:E:49:GLU:OE1	1:E:84:LEU:HD11	2.11	0.51
1:C:122:ARG:HH22	2:H:1:SGN:C2	2.23	0.51
1:F:86:LEU:HB2	1:F:98:ILE:HG23	1.92	0.51
1:A:98:ILE:HB	1:A:108:PHE:CE2	2.46	0.51
1:B:98:ILE:HG13	1:B:107:TRP:O	2.11	0.51
1:B:67:MSE:HG3	1:B:73:LEU:HD12	1.93	0.51
1:B:91:GLU:C	1:B:93:HIS:H	2.19	0.51
1:A:64:TYR:CZ	1:A:79:PRO:HG3	2.46	0.50
1:A:15:TYR:HB2	1:A:22:PHE:CE1	2.45	0.50
1:F:112:LYS:CE	1:F:116:SER:HB2	2.41	0.50
1:C:91:GLU:O	1:C:93:HIS:N	2.44	0.50
1:B:80:ASN:HD22	1:B:82:GLU:H	1.59	0.50
1:C:112:LYS:HG2	2:H:4:IDS:O1S	2.12	0.50
1:D:22:PHE:O	1:D:33:GLY:HA2	2.11	0.50
1:E:119:ARG:HD2	1:E:122:ARG:NH2	2.26	0.50
1:D:12:LYS:CE	1:D:46:LEU:HD22	2.42	0.50
1:A:112:LYS:HD2	1:A:116:SER:HB2	1.93	0.50
1:B:101:LYS:HE3	1:B:102:HIS:CE1	2.47	0.50
2:G:4:IDS:C6	2:G:5:SGN:H5	2.42	0.49
1:E:80:ASN:H	1:E:80:ASN:HD22	1.59	0.49
1:C:54:VAL:CG1	1:C:55:TYR:N	2.75	0.49
1:D:87:GLU:HG3	1:D:97:TYR:CE2	2.47	0.49
1:E:90:GLU:HG2	1:E:131:LEU:HD21	1.95	0.49
1:B:24:ARG:CZ	1:B:26:LEU:HD21	2.42	0.49
1:C:113:LYS:HG3	2:H:4:IDS:O2S	2.12	0.49
1:C:19:GLY:N	4:C:653:HOH:O	2.40	0.48
1:B:55:TYR:CE1	1:B:84:LEU:HD13	2.49	0.48
1:C:122:ARG:NH2	2:H:1:SGN:O1S	2.46	0.48
1:D:94:TYR:HA	1:D:132:PHE:O	2.13	0.48
1:D:89:LEU:HG	1:D:91:GLU:O	2.13	0.48
1:C:111:LEU:HD23	1:C:117:CYS:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:13:LEU:HB2	1:D:135:LEU:O	2.13	0.48
1:C:47:SER:HB3	1:C:55:TYR:HB2	1.96	0.48
1:A:95:ASN:HD22	1:A:134:PRO:CG	2.27	0.47
1:B:31:VAL:HG11	1:B:73:LEU:HD23	1.95	0.47
1:E:22:PHE:O	1:E:33:GLY:HA2	2.13	0.47
1:F:21:HIS:CE1	1:F:35:ARG:HE	2.32	0.47
1:A:47:SER:HB3	1:A:55:TYR:HB2	1.96	0.47
1:B:24:ARG:NE	1:B:26:LEU:HD21	2.30	0.47
1:F:54:VAL:HG22	1:F:85:PHE:O	2.15	0.47
1:C:46:LEU:HD22	1:C:56:ILE:HG12	1.96	0.47
1:E:73:LEU:HD11	1:E:109:VAL:HG13	1.94	0.47
1:D:112:LYS:CD	1:D:116:SER:HB2	2.43	0.47
1:E:26:LEU:N	1:E:26:LEU:HD22	2.28	0.47
1:E:17:SER:HB3	1:E:94:TYR:HE2	1.79	0.47
1:B:124:HIS:O	1:B:127:GLN:HG3	2.14	0.47
1:C:22:PHE:O	1:C:33:GLY:HA2	2.15	0.47
1:D:97:TYR:HB2	1:D:109:VAL:CG2	2.45	0.47
1:D:113:LYS:HZ3	3:I:5:IDS:H2	1.79	0.47
1:F:74:TYR:CE1	1:F:76:SER:HB2	2.47	0.47
1:A:11:PRO:HB3	1:A:59:THR:HG21	1.97	0.47
1:A:98:ILE:HG13	1:A:107:TRP:O	2.14	0.46
1:C:91:GLU:HB2	1:C:94:TYR:CD2	2.49	0.46
1:F:112:LYS:HE2	1:F:116:SER:HB2	1.97	0.46
1:A:84:LEU:HD23	1:A:100:LYS:HD2	1.97	0.46
3:I:5:IDS:H3	3:I:5:IDS:O1S	2.16	0.46
1:D:44:LEU:HD22	1:D:56:ILE:HG22	1.96	0.46
1:D:90:GLU:HG3	1:D:96:THR:HG21	1.97	0.46
1:C:24:ARG:HD3	1:C:26:LEU:CD2	2.46	0.46
1:A:119:ARG:O	1:A:120:GLY:C	2.58	0.46
1:E:89:LEU:HD12	1:E:94:TYR:O	2.16	0.46
1:F:47:SER:HB3	1:F:55:TYR:HB2	1.96	0.45
1:E:86:LEU:HD13	1:E:88:ARG:NH2	2.31	0.45
1:C:119:ARG:O	1:C:120:GLY:C	2.59	0.45
1:E:26:LEU:HB3	1:E:27:PRO:HD2	1.99	0.45
1:E:11:PRO:HD3	1:E:45:GLN:HE22	1.81	0.45
1:A:107:TRP:NE1	1:A:121:PRO:HA	2.31	0.45
1:B:118:LYS:HG2	1:B:123:THR:HG22	1.99	0.45
1:D:26:LEU:HB3	1:D:27:PRO:HD2	1.99	0.45
1:A:23:LEU:HD12	1:A:24:ARG:H	1.82	0.44
1:A:122:ARG:HD3	1:A:122:ARG:HA	1.64	0.44
1:B:89:LEU:HD23	1:B:95:ASN:OD1	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:107:TRP:HE1	1:D:121:PRO:HA	1.81	0.44
1:E:26:LEU:HD23	1:E:30:THR:HG22	1.99	0.44
1:B:14:LEU:O	1:B:22:PHE:HA	2.18	0.44
1:E:26:LEU:HD23	1:E:30:THR:O	2.18	0.44
1:E:106:ASN:HD21	1:E:125:TYR:N	2.15	0.44
1:A:18:ASN:HD22	1:A:129:ALA:HB2	1.82	0.44
1:C:24:ARG:HD3	1:C:26:LEU:HD21	1.99	0.44
1:C:25:ILE:HD11	1:C:65:LEU:HB2	1.99	0.44
1:C:66:ALA:O	1:C:73:LEU:HA	2.17	0.44
1:C:118:LYS:HE2	1:C:122:ARG:O	2.17	0.44
1:D:11:PRO:O	1:D:137:VAL:HG12	2.18	0.44
1:D:13:LEU:HD23	1:D:42:ILE:O	2.17	0.44
1:E:98:ILE:HD11	1:E:103:ALA:CA	2.47	0.44
1:F:27:PRO:HD3	1:F:41:HIS:CE1	2.52	0.44
1:B:18:ASN:HA	1:B:128:LYS:HD3	1.99	0.44
1:B:91:GLU:HA	1:B:91:GLU:OE1	2.17	0.44
1:F:80:ASN:ND2	1:F:83:CYS:SG	2.90	0.44
1:F:114:ASN:OD1	1:F:114:ASN:C	2.61	0.44
1:B:112:LYS:NZ	1:B:114:ASN:HD21	2.15	0.44
1:C:13:LEU:HD21	1:C:42:ILE:HG13	2.00	0.44
1:F:86:LEU:HB3	1:F:88:ARG:NH2	2.32	0.44
1:E:80:ASN:HD22	1:E:80:ASN:N	2.16	0.44
1:F:128:LYS:HA	1:F:131:LEU:CD1	2.48	0.44
1:C:80:ASN:HD21	1:C:82:GLU:HB2	1.83	0.44
1:E:21:HIS:HD2	1:E:35:ARG:HG2	1.83	0.44
1:C:91:GLU:CB	1:C:94:TYR:HD2	2.30	0.43
1:A:53:GLU:C	1:A:54:VAL:HG13	2.44	0.43
1:B:80:ASN:HD22	1:B:82:GLU:N	2.16	0.43
1:E:68:ASP:OD2	1:E:72:LEU:HB3	2.18	0.43
1:A:14:LEU:HD12	1:A:23:LEU:HD23	2.00	0.43
1:B:81:GLU:HG3	1:B:101:LYS:HD3	2.00	0.43
1:F:37:ARG:HG2	1:F:37:ARG:NH1	2.33	0.43
1:D:73:LEU:HD11	1:D:109:VAL:CG1	2.48	0.43
1:D:91:GLU:OE1	1:D:91:GLU:HA	2.17	0.43
1:D:97:TYR:HB2	1:D:109:VAL:HG23	2.00	0.43
1:E:129:ALA:HB2	2:H:2:IDS:O1S	2.18	0.43
2:H:2:IDS:O6A	2:H:3:SGN:H5	2.18	0.43
1:E:72:LEU:HB2	1:E:119:ARG:HH21	1.84	0.43
1:C:24:ARG:HH11	1:C:26:LEU:CD1	2.24	0.43
1:E:120:GLY:N	1:E:121:PRO:HD2	2.33	0.43
1:A:67:MSE:HE3	1:A:73:LEU:CD2	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:37:ARG:NH2	1:D:137:VAL:HA	2.33	0.43
1:D:67:MSE:HA	1:D:72:LEU:O	2.18	0.43
1:F:85:PHE:CD1	1:F:99:SER:HA	2.53	0.43
1:C:114:ASN:ND2	2:H:3:SGN:S2	2.92	0.42
1:F:46:LEU:HA	1:F:55:TYR:O	2.18	0.42
1:C:31:VAL:HG21	1:C:65:LEU:HD23	2.00	0.42
1:F:15:TYR:HB3	1:F:135:LEU:HD13	2.00	0.42
1:F:67:MSE:SE	1:F:107:TRP:HE3	2.51	0.42
3:I:1:IDS:O6A	3:I:2:SGN:H62	2.20	0.42
1:D:13:LEU:CD2	1:D:42:ILE:HG13	2.47	0.42
1:D:53:GLU:HB3	1:D:84:LEU:HD23	2.01	0.42
1:C:100:LYS:O	1:C:101:LYS:C	2.63	0.42
1:D:37:ARG:O	1:D:42:ILE:HD11	2.19	0.42
1:D:37:ARG:HH21	1:D:137:VAL:HG23	1.84	0.42
1:E:64:TYR:CE2	1:E:79:PRO:HD3	2.55	0.42
1:F:108:PHE:CD2	1:F:130:ILE:HD12	2.55	0.42
1:A:10:LYS:O	1:A:11:PRO:C	2.62	0.42
1:B:127:GLN:O	1:B:130:ILE:HG12	2.19	0.42
1:A:95:ASN:ND2	1:A:134:PRO:HG2	2.35	0.42
1:F:97:TYR:N	1:F:97:TYR:CD1	2.87	0.42
1:A:80:ASN:ND2	1:A:82:GLU:HB2	2.35	0.41
1:E:10:LYS:HE2	1:E:10:LYS:HB3	1.82	0.41
1:B:102:HIS:O	1:B:103:ALA:C	2.62	0.41
1:C:55:TYR:HE2	1:C:79:PRO:HB3	1.82	0.41
1:D:49:GLU:CB	1:D:55:TYR:HE1	2.32	0.41
1:E:79:PRO:HA	1:E:83:CYS:SG	2.60	0.41
1:E:86:LEU:HB2	1:E:98:ILE:HG23	2.01	0.41
1:C:24:ARG:HH21	1:C:34:THR:HB	1.85	0.41
1:C:27:PRO:C	1:C:29:GLY:H	2.28	0.41
1:D:96:THR:HG22	1:D:130:ILE:O	2.21	0.41
1:E:10:LYS:N	1:E:11:PRO:HD2	2.36	0.41
1:B:12:LYS:O	1:B:13:LEU:HD12	2.21	0.41
1:B:86:LEU:HD21	1:B:100:LYS:HG3	2.02	0.41
1:D:127:GLN:O	1:D:130:ILE:HG12	2.20	0.41
1:F:43:GLN:OE1	1:F:137:VAL:HG11	2.20	0.41
1:B:90:GLU:OE1	1:B:94:TYR:HD2	2.04	0.41
1:D:47:SER:HB3	1:D:55:TYR:HB2	2.02	0.41
1:F:15:TYR:CE1	1:F:20:GLY:HA2	2.56	0.41
1:D:66:ALA:O	1:D:73:LEU:HA	2.21	0.41
1:A:107:TRP:HE1	1:A:121:PRO:HA	1.85	0.41
1:B:44:LEU:HD23	1:B:56:ILE:CG2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:111:LEU:HD23	1:F:117:CYS:HA	2.03	0.41
1:B:46:LEU:N	1:B:46:LEU:HD12	2.35	0.41
1:D:24:ARG:O	1:D:31:VAL:HA	2.20	0.41
1:D:24:ARG:HD3	1:D:26:LEU:CD2	2.51	0.41
1:D:24:ARG:HD3	1:D:26:LEU:HD21	2.01	0.41
1:B:98:ILE:HA	1:B:108:PHE:HA	2.03	0.41
1:A:73:LEU:HG	4:A:622:HOH:O	2.20	0.40
1:F:24:ARG:CZ	1:F:26:LEU:HD21	2.51	0.40
1:B:111:LEU:HD23	1:B:111:LEU:HA	1.85	0.40
1:C:26:LEU:HB3	1:C:27:PRO:HD2	2.03	0.40
1:C:89:LEU:HA	1:C:94:TYR:O	2.21	0.40
1:C:124:HIS:O	1:C:127:GLN:HG3	2.21	0.40
1:F:14:LEU:HD12	1:F:23:LEU:HD23	2.02	0.40
1:F:124:HIS:ND1	1:F:125:TYR:N	2.68	0.40
2:G:4:IDS:O6B	2:G:5:SGN:H5	2.21	0.40
1:B:90:GLU:OE1	1:B:94:TYR:HB2	2.20	0.40
1:C:50:SER:C	1:C:52:GLY:N	2.79	0.40
1:C:109:VAL:HA	4:C:648:HOH:O	2.22	0.40
1:D:106:ASN:ND2	1:D:106:ASN:O	2.54	0.40
1:E:16:CYS:HB2	1:E:132:PHE:CE1	2.56	0.40
2:H:2:IDS:C6	2:H:3:SGN:H5	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	128/135 (95%)	117 (91%)	11 (9%)	0	100	100
1	B	125/135 (93%)	112 (90%)	12 (10%)	1 (1%)	16	50
1	C	126/135 (93%)	106 (84%)	15 (12%)	5 (4%)	2	14
1	D	125/135 (93%)	109 (87%)	14 (11%)	2 (2%)	7	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	127/135 (94%)	116 (91%)	9 (7%)	2 (2%)	7	34
1	F	127/135 (94%)	113 (89%)	13 (10%)	1 (1%)	16	50
All	All	758/810 (94%)	673 (89%)	74 (10%)	11 (2%)	8	35

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	92	ASN
1	B	36	ASP
1	D	36	ASP
1	E	90	GLU
1	F	50	SER
1	C	68	ASP
1	D	68	ASP
1	E	36	ASP
1	C	51	VAL
1	C	52	GLY
1	C	120	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	112/117 (96%)	109 (97%)	3 (3%)	39	71
1	B	110/117 (94%)	102 (93%)	8 (7%)	13	42
1	C	111/117 (95%)	100 (90%)	11 (10%)	7	30
1	D	110/117 (94%)	105 (96%)	5 (4%)	24	59
1	E	113/117 (97%)	105 (93%)	8 (7%)	13	43
1	F	113/117 (97%)	110 (97%)	3 (3%)	39	71
All	All	669/702 (95%)	631 (94%)	38 (6%)	18	52

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
1	A	135	LEU
1	A	137	VAL
1	B	37	ARG
1	B	65	LEU
1	B	70	ASP
1	B	78	THR
1	B	80	ASN
1	B	90	GLU
1	B	112	LYS
1	B	113	LYS
1	C	46	LEU
1	C	47	SER
1	C	57	LYS
1	C	80	ASN
1	C	89	LEU
1	C	91	GLU
1	C	95	ASN
1	C	98	ILE
1	C	106	ASN
1	C	116	SER
1	C	128	LYS
1	D	45	GLN
1	D	49	GLU
1	D	57	LYS
1	D	84	LEU
1	D	109	VAL
1	E	30	THR
1	E	60	GLU
1	E	65	LEU
1	E	80	ASN
1	E	98	ILE
1	E	119	ARG
1	E	130	ILE
1	E	133	LEU
1	F	13	LEU
1	F	80	ASN
1	F	92	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	HIS

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Mol	Chain	Res	Type
1	A	43	GLN
1	A	80	ASN
1	A	95	ASN
1	A	102	HIS
1	A	124	HIS
1	B	21	HIS
1	B	80	ASN
1	B	93	HIS
1	B	114	ASN
1	C	41	HIS
1	C	80	ASN
1	C	95	ASN
1	C	102	HIS
1	C	106	ASN
1	D	18	ASN
1	D	41	HIS
1	D	45	GLN
1	D	106	ASN
1	D	124	HIS
1	E	80	ASN
1	E	92	ASN
1	E	93	HIS
1	E	95	ASN
1	E	106	ASN
1	F	41	HIS
1	F	63	GLN
1	F	80	ASN
1	F	95	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

15 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SGN	G	1	2	20,20,20	1.83	4 (20%)	25,31,31	1.47	4 (16%)
2	IDS	G	2	2	16,16,17	1.60	4 (25%)	16,24,26	1.14	1 (6%)
2	SGN	G	3	2	19,19,20	1.16	3 (15%)	23,29,31	1.29	2 (8%)
2	IDS	G	4	2	16,16,17	1.86	3 (18%)	16,24,26	1.12	2 (12%)
2	SGN	G	5	2	19,19,20	1.11	1 (5%)	23,29,31	1.44	2 (8%)
2	SGN	H	1	2	20,20,20	2.18	3 (15%)	25,31,31	1.21	2 (8%)
2	IDS	H	2	2	16,16,17	1.55	2 (12%)	16,24,26	1.07	1 (6%)
2	SGN	H	3	2	19,19,20	1.29	2 (10%)	23,29,31	1.45	3 (13%)
2	IDS	H	4	2	16,16,17	1.31	3 (18%)	16,24,26	1.30	4 (25%)
2	SGN	H	5	2	19,19,20	1.12	2 (10%)	23,29,31	1.45	5 (21%)
3	IDS	I	1	3	17,17,17	1.04	2 (11%)	22,26,26	1.29	3 (13%)
3	SGN	I	2	3	19,19,20	1.63	3 (15%)	23,29,31	1.20	1 (4%)
3	IDS	I	3	3	16,16,17	1.21	3 (18%)	16,24,26	1.34	3 (18%)
3	SGN	I	4	3	19,19,20	2.07	3 (15%)	23,29,31	1.71	5 (21%)
3	IDS	I	5	3	16,16,17	1.43	2 (12%)	16,24,26	1.23	2 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SGN	G	1	2	-	4/11/31/31	0/1/1/1
2	IDS	G	2	2	-	5/9/26/29	0/1/1/1
2	SGN	G	3	2	-	1/11/28/31	0/1/1/1
2	IDS	G	4	2	-	0/9/26/29	0/1/1/1
2	SGN	G	5	2	-	3/11/28/31	0/1/1/1
2	SGN	H	1	2	-	3/11/31/31	0/1/1/1
2	IDS	H	2	2	-	3/9/26/29	0/1/1/1
2	SGN	H	3	2	-	6/11/28/31	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IDS	H	4	2	-	0/9/26/29	0/1/1/1
2	SGN	H	5	2	-	2/11/28/31	0/1/1/1
3	IDS	I	1	3	-	3/9/29/29	0/1/1/1
3	SGN	I	2	3	-	3/11/28/31	0/1/1/1
3	IDS	I	3	3	-	0/9/26/29	0/1/1/1
3	SGN	I	4	3	-	4/11/28/31	0/1/1/1
3	IDS	I	5	3	1/1/6/7	2/9/26/29	0/1/1/1

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	1	SGN	S1-N2	7.79	1.69	1.59
3	I	4	SGN	S1-N2	6.79	1.68	1.59
2	G	1	SGN	O1S-S1	4.88	1.47	1.42
2	H	2	IDS	O2-C2	-4.50	1.40	1.47
2	G	4	IDS	O2-C2	-4.34	1.40	1.47
3	I	2	SGN	O2S-S1	4.03	1.46	1.42
3	I	5	IDS	O2-C2	-3.63	1.41	1.47
2	G	2	IDS	O2-C2	-3.55	1.42	1.47
2	G	1	SGN	S1-N2	3.53	1.63	1.59
2	G	4	IDS	O6B-C6	-3.41	1.19	1.30
2	G	4	IDS	C5-C6	-3.16	1.46	1.53
2	G	1	SGN	O2S-S1	3.09	1.45	1.42
2	G	5	SGN	S1-N2	3.08	1.63	1.59
2	H	1	SGN	O1S-S1	3.02	1.45	1.42
3	I	2	SGN	O1S-S1	3.01	1.45	1.42
2	H	3	SGN	O2S-S1	3.00	1.45	1.42
3	I	2	SGN	S1-N2	2.96	1.63	1.59
3	I	3	IDS	O2-C2	-2.84	1.43	1.47
2	H	1	SGN	C2-N2	2.81	1.51	1.47
3	I	4	SGN	O1S-S1	2.76	1.45	1.42
3	I	1	IDS	O6B-C6	-2.71	1.22	1.30
2	H	4	IDS	C1-C2	2.64	1.55	1.51
3	I	5	IDS	O6B-C6	-2.61	1.22	1.30
2	H	4	IDS	O6B-C6	-2.60	1.22	1.30
2	G	2	IDS	O6B-C6	-2.57	1.22	1.30
2	H	2	IDS	O6B-C6	-2.54	1.22	1.30
3	I	1	IDS	O2-C2	-2.53	1.40	1.46
2	G	1	SGN	O6-C6	-2.50	1.37	1.46
3	I	4	SGN	O2S-S1	2.44	1.44	1.42
3	I	3	IDS	C3-C2	2.41	1.58	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	4	IDS	O2-C2	-2.39	1.43	1.47
2	G	3	SGN	O1S-S1	2.35	1.44	1.42
2	H	5	SGN	O2S-S1	2.31	1.44	1.42
2	G	3	SGN	O6-C6	-2.19	1.38	1.46
2	H	3	SGN	O1S-S1	2.18	1.44	1.42
3	I	3	IDS	O6B-C6	-2.17	1.23	1.30
2	G	3	SGN	C4-C3	2.13	1.57	1.52
2	H	5	SGN	O1S-S1	2.10	1.44	1.42
2	G	2	IDS	C4-C5	-2.08	1.49	1.53
2	G	2	IDS	C5-C6	-2.04	1.48	1.53

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	5	SGN	O1S-S1-O2S	-4.72	109.86	120.36
2	G	1	SGN	C4-C3-C2	-4.17	104.33	110.40
2	G	3	SGN	O1S-S1-O2S	-4.11	111.22	120.36
2	H	1	SGN	O1S-S1-O2S	-3.88	111.74	120.36
3	I	4	SGN	C3-C2-N2	3.82	115.34	110.32
3	I	2	SGN	O1S-S1-O2S	-3.74	112.05	120.36
3	I	4	SGN	C4-C3-C2	-3.65	105.67	111.02
3	I	1	IDS	C2-O2-S	3.63	127.75	119.04
3	I	4	SGN	O1S-S1-O2S	-3.56	112.44	120.36
2	H	3	SGN	O1S-S1-O2S	-3.47	112.64	120.36
2	H	5	SGN	O1S-S1-O2S	-3.42	112.75	120.36
2	G	1	SGN	C3-C4-C5	-3.03	104.74	110.23
2	G	1	SGN	O1S-S1-O2S	-2.86	113.99	120.36
2	H	3	SGN	C1-O5-C5	2.86	116.01	112.19
3	I	1	IDS	O6B-C6-O6A	-2.80	117.72	124.08
3	I	3	IDS	O6B-C6-O6A	-2.71	117.94	124.08
3	I	5	IDS	O6B-C6-O6A	-2.69	117.97	124.08
3	I	3	IDS	O2-C2-C3	2.51	110.46	106.95
2	H	2	IDS	O6B-C6-O6A	-2.49	118.44	124.08
2	H	5	SGN	C3-C4-C5	-2.45	105.79	110.23
2	G	3	SGN	O2S-S1-N2	-2.41	104.85	108.88
2	H	5	SGN	C1-O5-C5	2.37	115.36	112.19
3	I	4	SGN	O2S-S1-N2	-2.37	104.92	108.88
2	G	5	SGN	C4-C3-C2	-2.36	107.56	111.02
3	I	5	IDS	O6B-C6-C5	2.32	122.02	113.64
2	H	4	IDS	O6B-C6-O6A	-2.32	118.82	124.08
2	H	3	SGN	O2S-S1-N2	-2.21	105.19	108.88
3	I	4	SGN	O4-C4-C3	2.20	115.56	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1	SGN	C1-C2-C3	-2.19	107.56	110.54
2	G	4	IDS	O3S-S-O2	2.17	111.36	106.37
2	H	5	SGN	C4-C3-C2	-2.15	107.86	111.02
2	G	4	IDS	O6B-C6-O6A	-2.14	119.22	124.08
2	H	4	IDS	C3-C4-C5	2.13	112.97	109.30
2	H	4	IDS	O6B-C6-C5	2.13	121.33	113.64
2	G	1	SGN	O2S-S1-N2	-2.13	105.32	108.88
2	G	2	IDS	O6B-C6-O6A	-2.09	119.34	124.08
2	H	5	SGN	O6-C6-C5	2.09	111.29	107.57
2	H	4	IDS	C4-C3-C2	2.08	114.01	110.23
3	I	3	IDS	O6B-C6-C5	2.07	121.11	113.64
3	I	1	IDS	O6B-C6-C5	2.04	121.00	113.64

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	I	5	IDS	C1

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	1	SGN	C6-O6-S2-O4S
2	G	1	SGN	C6-O6-S2-O6S
2	G	2	IDS	C1-C2-O2-S
2	G	2	IDS	C3-C2-O2-S
2	G	2	IDS	C2-O2-S-O1S
2	G	2	IDS	C2-O2-S-O2S
2	H	1	SGN	C2-N2-S1-O2S
2	H	1	SGN	C2-N2-S1-O3S
2	H	3	SGN	C4-C5-C6-O6
2	H	3	SGN	O5-C5-C6-O6
2	H	3	SGN	C2-N2-S1-O2S
2	H	5	SGN	C2-N2-S1-O1S
3	I	1	IDS	C3-C2-O2-S
3	I	4	SGN	C6-O6-S2-O4S
3	I	4	SGN	C6-O6-S2-O6S
3	I	5	IDS	C1-C2-O2-S
2	G	1	SGN	C6-O6-S2-O5S
3	I	4	SGN	C6-O6-S2-O5S
2	H	2	IDS	C2-O2-S-O2S
2	G	5	SGN	C6-O6-S2-O6S
2	H	3	SGN	C6-O6-S2-O6S

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Mol	Chain	Res	Type	Atoms
2	H	3	SGN	C6-O6-S2-O5S
2	G	2	IDS	C2-O2-S-O3S
2	H	2	IDS	C2-O2-S-O1S
2	H	2	IDS	C2-O2-S-O3S
2	H	3	SGN	C6-O6-S2-O4S
2	G	5	SGN	C6-O6-S2-O5S
3	I	5	IDS	C3-C2-O2-S
3	I	1	IDS	C1-C2-O2-S
2	G	5	SGN	C6-O6-S2-O4S
2	H	1	SGN	C2-N2-S1-O1S
2	H	5	SGN	C4-C5-C6-O6
3	I	1	IDS	C2-O2-S-O3S
3	I	2	SGN	C6-O6-S2-O5S
3	I	2	SGN	C6-O6-S2-O4S
3	I	2	SGN	C2-N2-S1-O1S
3	I	4	SGN	C2-N2-S1-O2S
2	G	1	SGN	C2-N2-S1-O2S
2	G	3	SGN	C2-N2-S1-O2S

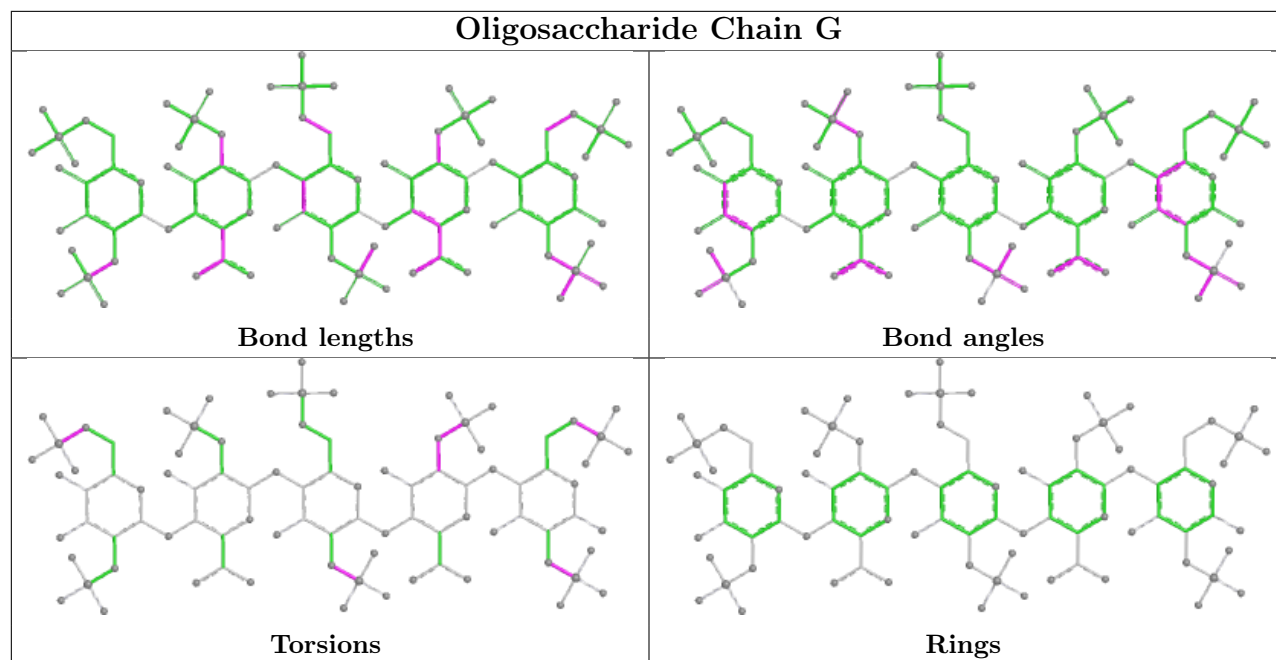
There are no ring outliers.

13 monomers are involved in 22 short contacts:

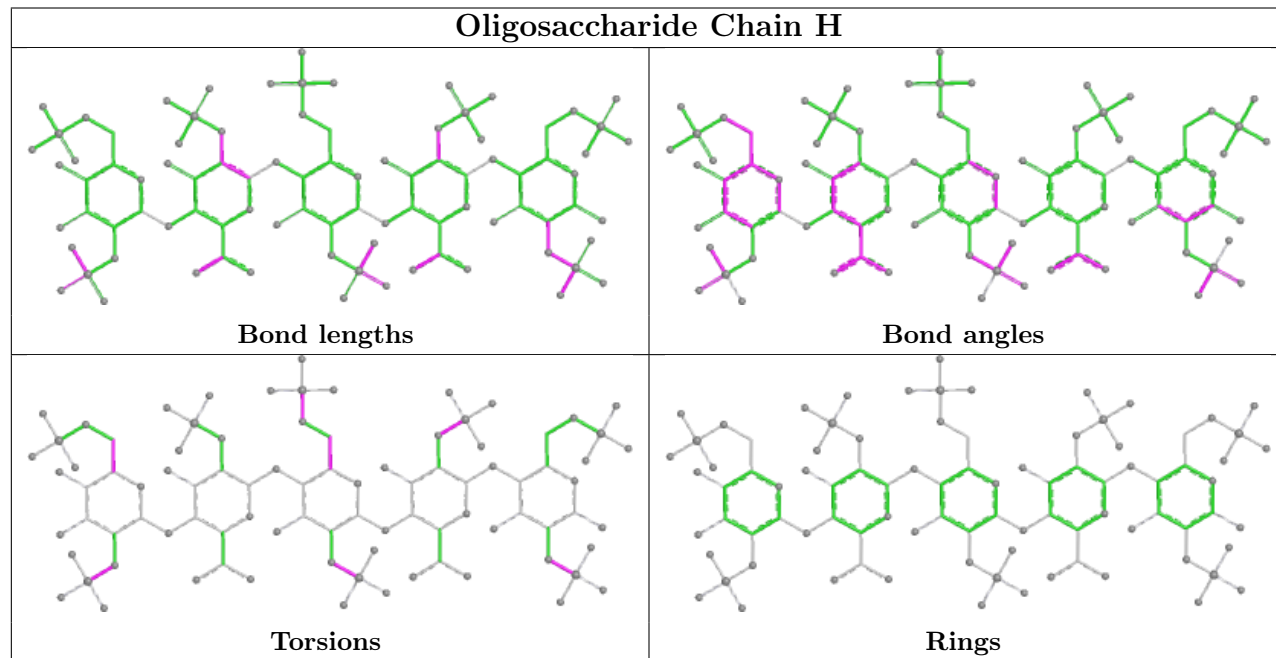
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	4	IDS	2	0
2	H	3	SGN	3	0
2	H	5	SGN	1	0
2	H	1	SGN	2	0
2	H	2	IDS	3	0
2	H	4	IDS	4	0
2	G	5	SGN	2	0
3	I	1	IDS	1	0
3	I	5	IDS	5	0
3	I	2	SGN	1	0
2	G	2	IDS	1	0
2	G	1	SGN	2	0
2	G	3	SGN	2	0

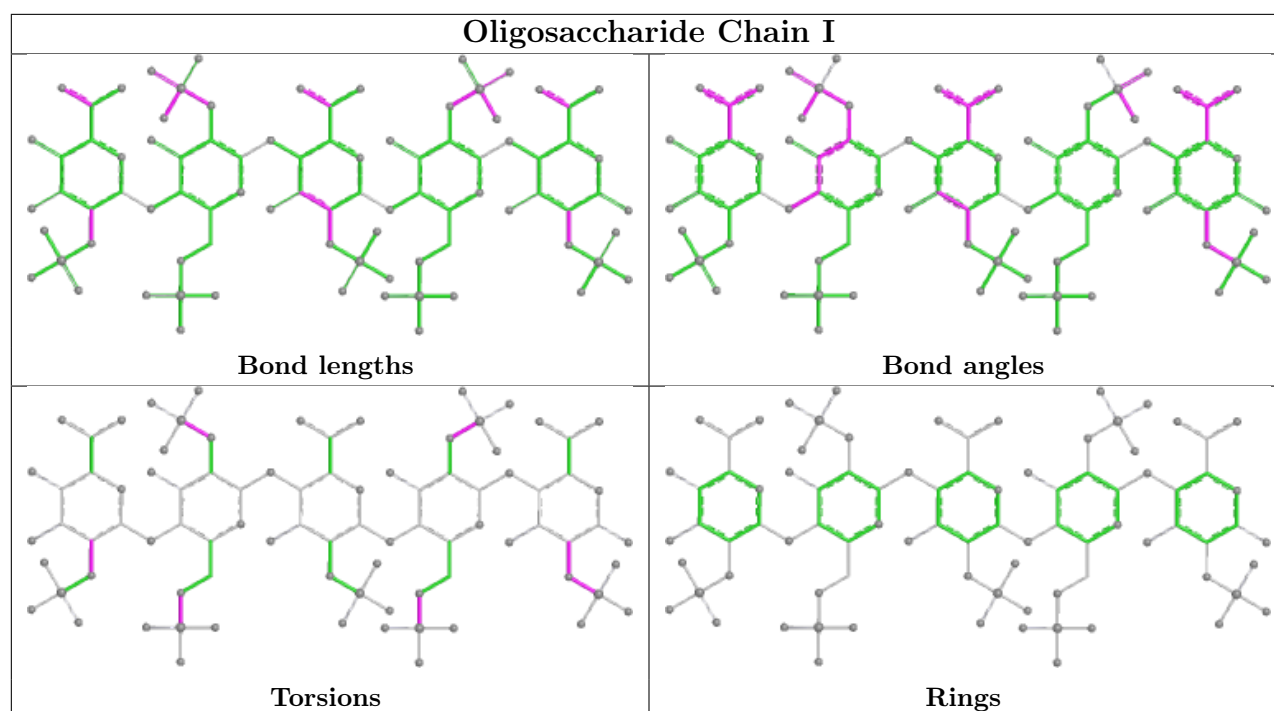
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain G



Oligosaccharide Chain H





5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.