



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:52 AM GMT

PDB ID : 3Q3G
Title : Crystal Structure of A-domain in complex with antibody
Authors : Mahalingam, B.; Xiong, J.P.; Arnaout, M.A.
Deposited on : 2010-12-21
Resolution : 2.70 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

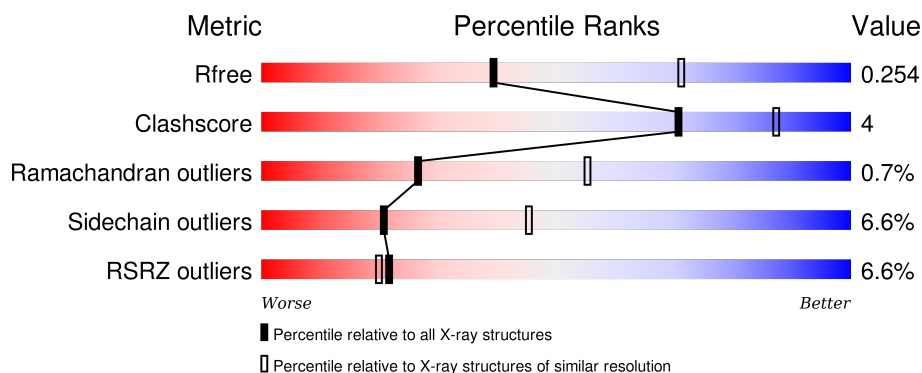
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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.




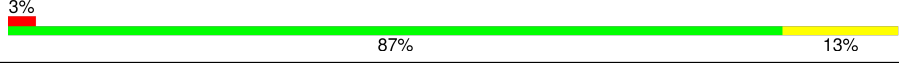

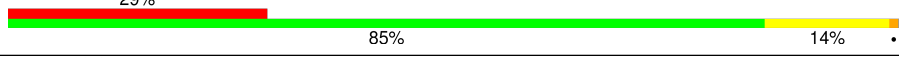
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	220	<div> <div></div> <div>82%16%•</div> </div>
1	C	220	<div> <div></div> <div>84%15%•</div> </div>
1	F	220	<div> <div>16%</div> <div>83%15%•</div> </div>
1	J	220	<div> <div></div> <div>81%16%•</div> </div>
2	B	224	<div> <div>4%</div> <div>85%13%•</div> </div>

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Mol	Chain	Length	Quality of chain	
2	D	224		•
2	H	224		•
2	K	224		
3	E	190		••
3	G	190		••
3	I	190		••
3	L	190		••

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
4	EDO	A	618	-	-	-	X
4	EDO	B	607	-	-	-	X
4	EDO	B	616	-	-	-	X
4	EDO	C	609	-	-	-	X
4	EDO	C	628	-	-	-	X
4	EDO	D	613	-	-	-	X
4	EDO	D	614	-	-	-	X
4	EDO	D	619	-	-	-	X
4	EDO	E	600	-	-	-	X
4	EDO	E	627	-	-	-	X
4	EDO	G	620	-	-	-	X
5	GOL	A	700	-	-	-	X
5	GOL	B	704	-	-	-	X
5	GOL	J	709	-	-	-	X
5	GOL	K	706	-	-	-	X
5	GOL	K	707	-	-	-	X

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 20257 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 Antibody Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	220	Total	C	N	O	S	0	0	0
			1718	1074	283	353	8			
1	A	220	Total	C	N	O	S	0	0	0
			1718	1074	283	353	8			
1	F	220	Total	C	N	O	S	0	0	0
			1718	1074	283	353	8			
1	J	220	Total	C	N	O	S	0	0	0
			1718	1074	283	353	8			

- Molecule 2 is a protein called Antibody Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	224	Total	C	N	O	S	0	0	0
			1689	1062	277	342	8			
2	B	224	Total	C	N	O	S	0	0	0
			1689	1062	277	342	8			
2	H	224	Total	C	N	O	S	0	0	0
			1689	1062	277	342	8			
2	K	224	Total	C	N	O	S	0	0	0
			1689	1062	277	342	8			

- Molecule 3 is a protein called Integrin alpha-M.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	189	Total	C	N	O	S	0	0	0
			1531	975	271	282	3			
3	E	189	Total	C	N	O	S	0	0	0
			1531	975	271	282	3			
3	I	189	Total	C	N	O	S	0	0	0
			1531	975	271	282	3			
3	L	189	Total	C	N	O	S	0	0	0
			1531	975	271	282	3			

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



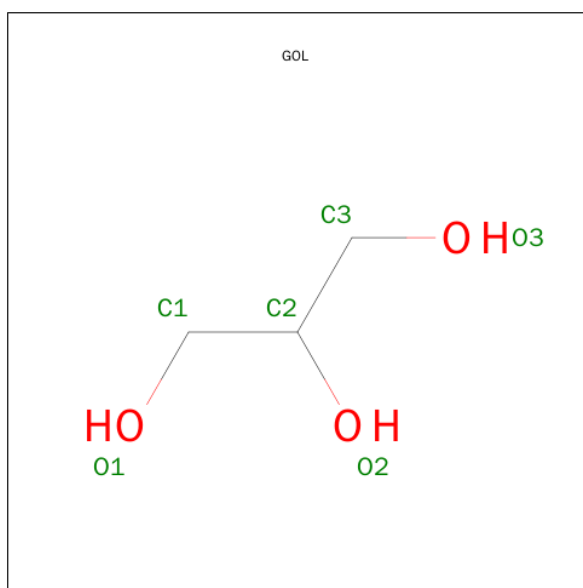
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	G	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		
4	J	1	Total	C	O	0	0
			4	2	2		
4	J	1	Total	C	O	0	0
			4	2	2		
4	K	1	Total	C	O	0	0
			4	2	2		
4	K	1	Total	C	O	0	0
			4	2	2		
4	K	1	Total	C	O	0	0
			4	2	2		
4	L	1	Total	C	O	0	0
			4	2	2		
4	L	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total C O 6 3 3	0	0
5	D	1	Total C O 6 3 3	0	0
5	D	1	Total C O 6 3 3	0	0
5	D	1	Total C O 6 3 3	0	0
5	G	1	Total C O 6 3 3	0	0
5	A	1	Total C O 6 3 3	0	0
5	B	1	Total C O 6 3 3	0	0
5	B	1	Total C O 6 3 3	0	0
5	J	1	Total C O 6 3 3	0	0
5	K	1	Total C O 6 3 3	0	0
5	K	1	Total C O 6 3 3	0	0

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	H	2	Total Na 2 2	0	0
6	D	1	Total Na 1 1	0	0

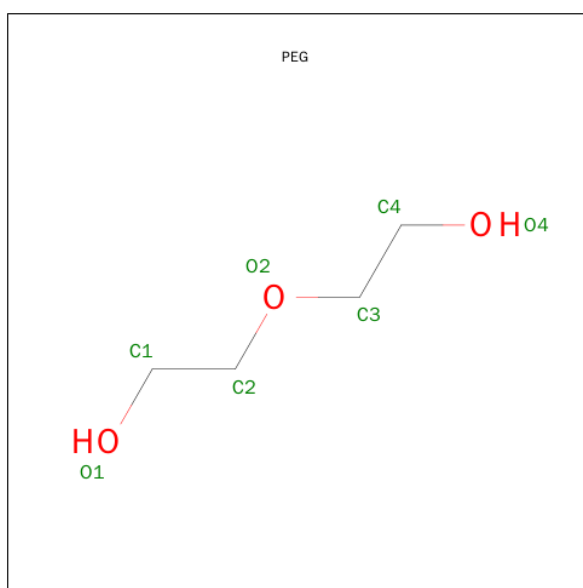
- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	G	1	Total Ca 1 1	0	0
7	I	1	Total Ca 1 1	0	0
7	L	1	Total Ca 1 1	0	0
7	E	1	Total Ca 1 1	0	0

- Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	G	1	Total Cl 1 1	0	0
8	J	1	Total Cl 1 1	0	0
8	K	1	Total Cl 1 1	0	0
8	E	2	Total Cl 2 2	0	0

- Molecule 9 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	H	1	Total C O 7 4 3	0	0

- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	G	29	Total O 29 29	0	0
10	E	12	Total O 12 12	0	0
10	I	6	Total O 6 6	0	0
10	L	10	Total O 10 10	0	0
10	C	42	Total O 42 42	0	0

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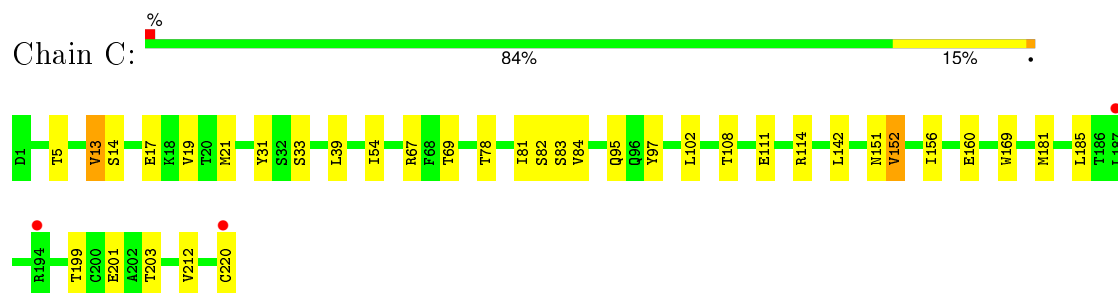
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	31	Total 31	O 31	0	0
10	F	6	Total 6	O 6	0	0
10	J	31	Total 31	O 31	0	0
10	D	42	Total 42	O 42	0	0
10	B	49	Total 49	O 49	0	0
10	H	22	Total 22	O 22	0	0
10	K	40	Total 40	O 40	0	0

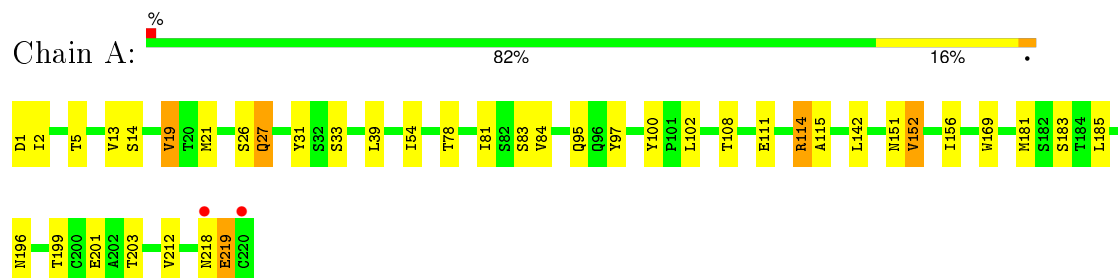
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

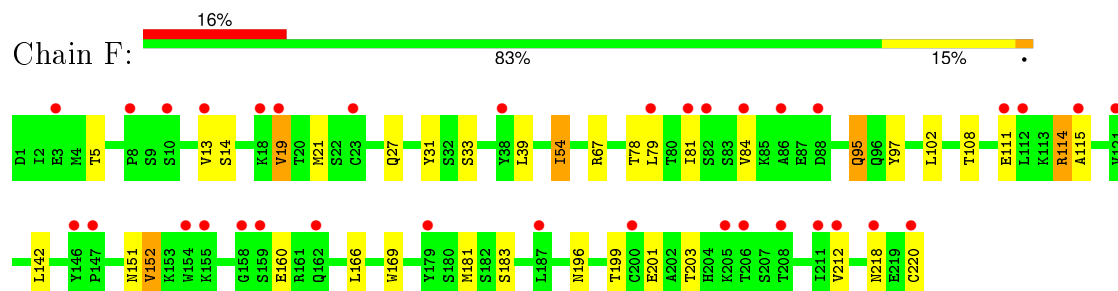
- Molecule 1: Antibody Light Chain



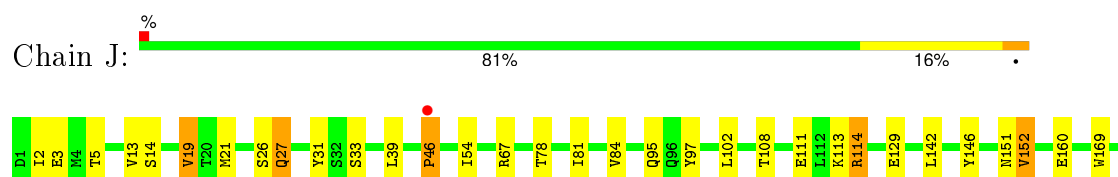
- Molecule 1: Antibody Light Chain

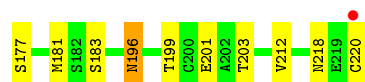


- Molecule 1: Antibody Light Chain

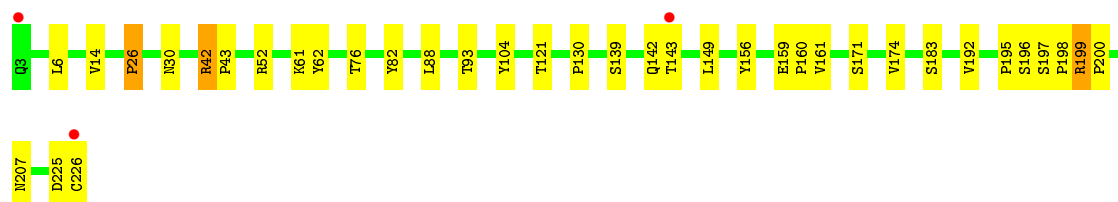
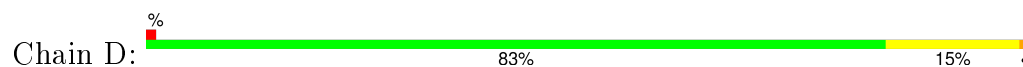


- Molecule 1: Antibody Light Chain

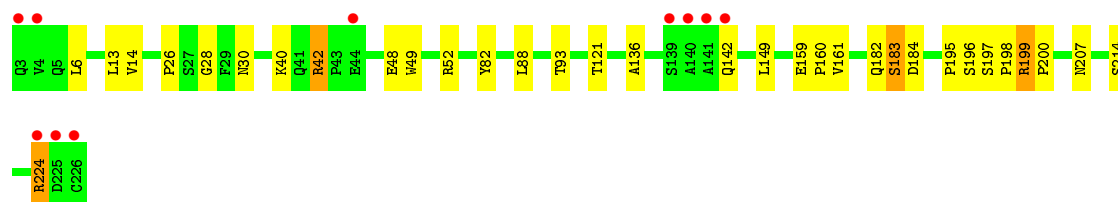
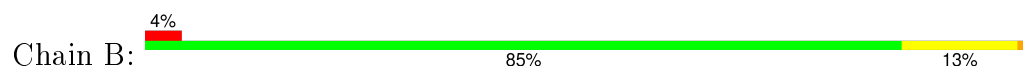




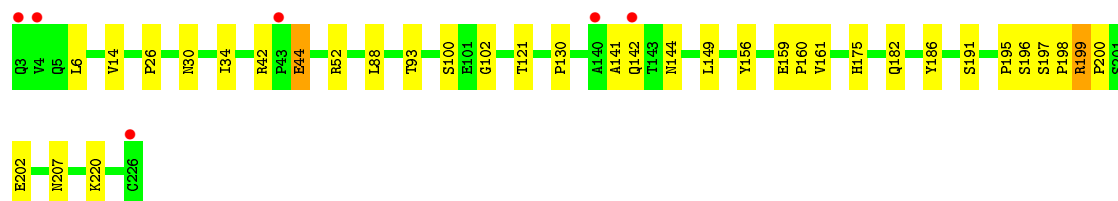
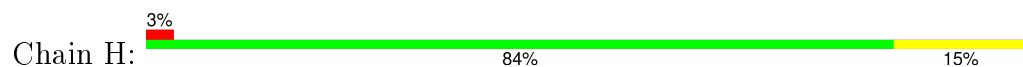
• Molecule 2: Antibody Heavy chain



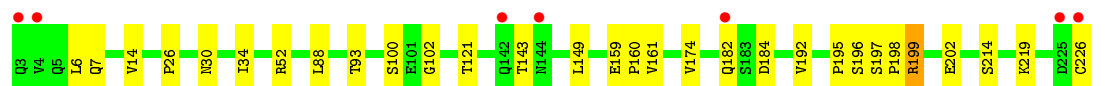
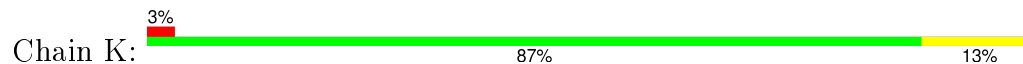
• Molecule 2: Antibody Heavy chain



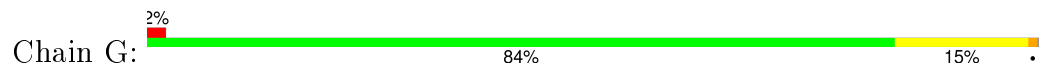
• Molecule 2: Antibody Heavy chain



• Molecule 2: Antibody Heavy chain

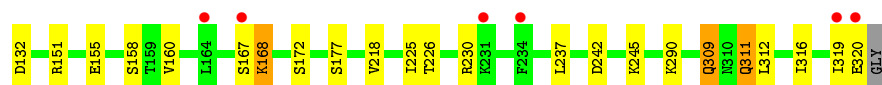


• Molecule 3: Integrin alpha-M



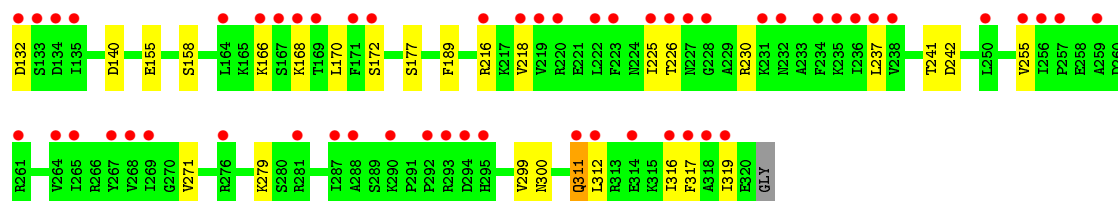
• Molecule 3: Integrin alpha-M

Chain E: 3% 87% 11% ..



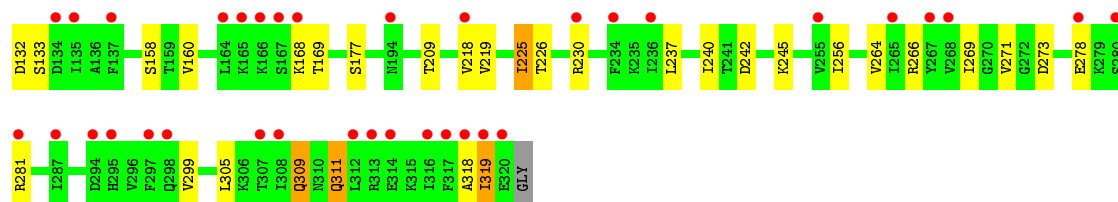
• Molecule 3: Integrin alpha-M

Chain I: 29% 85% 14% ..



• Molecule 3: Integrin alpha-M

Chain L: 18% 83% 14% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.94Å 157.22Å 232.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.76 – 2.70 47.76 – 2.70	Depositor EDS
% Data completeness (in resolution range)	(Not available) (47.76-2.70) 97.2 (47.76-2.70)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 2.69Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, R_{free}	0.211 , 0.244 0.221 , 0.254	Depositor DCC
R_{free} test set	4048 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	47.3	Xtriage
Anisotropy	0.223	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 80985 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	20257	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CL, NA, CA, EDO, PEG

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	0.44	0/1757	0.71	1/2384 (0.0%)
1	C	0.44	0/1757	0.69	0/2384
1	F	0.40	0/1757	0.68	0/2384
1	J	0.44	0/1757	0.74	0/2384
2	B	0.44	0/1732	0.74	0/2368
2	D	0.45	0/1732	0.75	0/2368
2	H	0.44	0/1732	0.74	0/2368
2	K	0.45	0/1732	0.74	0/2368
3	E	0.44	0/1560	0.71	1/2099 (0.0%)
3	G	0.46	0/1560	0.70	0/2099
3	I	0.45	0/1560	0.67	0/2099
3	L	0.47	0/1560	0.70	0/2099
All	All	0.44	0/20196	0.71	2/27404 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	219	GLU	C-N-CA	5.51	135.47	121.70
3	E	167	SER	N-CA-C	-5.27	96.76	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1718	0	1647	14	0
1	C	1718	0	1647	14	0
1	F	1718	0	1647	14	0
1	J	1718	0	1647	14	0
2	B	1689	0	1638	21	0
2	D	1689	0	1638	21	0
2	H	1689	0	1638	20	0
2	K	1689	0	1638	18	0
3	E	1531	0	1545	6	0
3	G	1531	0	1545	14	0
3	I	1531	0	1545	10	0
3	L	1531	0	1545	16	0
4	A	8	0	12	0	0
4	B	16	0	24	3	0
4	C	8	0	12	0	0
4	D	20	0	30	3	0
4	E	8	0	12	1	0
4	F	4	0	6	0	0
4	G	4	0	6	0	0
4	J	8	0	12	0	0
4	K	16	0	24	0	0
4	L	8	0	12	0	0
5	A	6	0	8	1	0
5	B	12	0	16	1	0
5	C	6	0	8	0	0
5	D	18	0	24	1	0
5	G	6	0	8	1	0
5	J	6	0	8	0	0
5	K	12	0	16	1	0
6	D	1	0	0	0	0
6	H	2	0	0	0	0
7	E	1	0	0	0	0
7	G	1	0	0	0	0
7	I	1	0	0	0	0
7	L	1	0	0	0	0
8	E	2	0	0	0	0
8	G	1	0	0	0	0
8	J	1	0	0	0	0
8	K	1	0	0	0	0
9	H	7	0	10	0	0
10	A	31	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	B	49	0	0	0	0
10	C	42	0	0	1	0
10	D	42	0	0	0	0
10	E	12	0	0	0	0
10	F	6	0	0	1	0
10	G	29	0	0	0	0
10	H	22	0	0	0	0
10	I	6	0	0	0	0
10	J	31	0	0	0	0
10	K	40	0	0	0	0
10	L	10	0	0	1	0
All	All	20257	0	19568	174	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 4.

All (174) 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)
2:B:199:ARG:HH11	2:B:199:ARG:HG3	1.00	1.12
2:D:199:ARG:HH11	2:D:199:ARG:HG3	1.09	1.10
1:C:13:VAL:HG22	1:C:19:VAL:HG11	1.41	1.02
2:K:199:ARG:HG3	2:K:199:ARG:HH11	1.23	1.00
2:H:199:ARG:HG2	2:H:200:PRO:HA	1.43	0.97
2:B:199:ARG:HH11	2:B:199:ARG:CG	1.86	0.89
2:D:199:ARG:HG3	2:D:199:ARG:NH1	1.89	0.87
2:B:199:ARG:NH1	2:B:199:ARG:HG3	1.80	0.87
1:C:13:VAL:CG2	1:C:19:VAL:HG11	2.05	0.86
2:K:199:ARG:CG	2:K:199:ARG:HH11	1.88	0.85
2:H:149:LEU:HD11	2:H:199:ARG:HD3	1.62	0.81
5:D:702:GOL:H12	2:H:202:GLU:OE1	1.83	0.78
2:D:199:ARG:CG	2:D:199:ARG:HH11	1.93	0.77
1:C:21:MET:HG2	1:C:108:THR:HG21	1.67	0.77
1:F:142:LEU:HD11	1:F:152:VAL:HG13	1.70	0.73
1:F:21:MET:HG2	1:F:108:THR:HG21	1.70	0.73
1:A:142:LEU:HD11	1:A:152:VAL:HG13	1.70	0.73
2:K:199:ARG:HG3	2:K:199:ARG:NH1	1.97	0.72
1:C:142:LEU:HD11	1:C:152:VAL:HG13	1.70	0.72
1:C:201:GLU:HG3	1:C:212:VAL:HG22	1.73	0.71
1:J:142:LEU:HD11	1:J:152:VAL:HG13	1.72	0.70
1:A:201:GLU:HG3	1:A:212:VAL:HG22	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:201:GLU:HG3	1:J:212:VAL:HG22	1.73	0.70
1:F:54:ILE:HD13	1:F:79:LEU:HD11	1.73	0.69
1:F:201:GLU:HG3	1:F:212:VAL:HG22	1.73	0.69
1:J:19:VAL:HG13	1:J:81:ILE:HB	1.76	0.67
2:B:195:PRO:HB2	2:B:198:PRO:HD2	1.77	0.66
2:H:34:ILE:HG22	2:H:102:GLY:HA2	1.76	0.66
2:H:195:PRO:HB2	2:H:198:PRO:HD2	1.77	0.66
2:D:195:PRO:HB2	2:D:198:PRO:HD2	1.78	0.66
1:J:21:MET:HG2	1:J:108:THR:HG21	1.78	0.65
1:A:100:TYR:HB2	4:E:600:EDO:H21	1.78	0.65
2:K:195:PRO:HB2	2:K:198:PRO:HD2	1.80	0.64
3:L:271:VAL:HG13	3:L:299:VAL:HG23	1.83	0.61
2:H:42:ARG:O	2:H:44:GLU:HA	2.00	0.61
2:D:149:LEU:HD11	2:D:199:ARG:HG2	1.82	0.61
1:F:169:TRP:CD1	1:F:181:MET:HG3	2.36	0.61
3:G:218:VAL:HG11	3:G:237:LEU:HD13	1.83	0.60
1:F:166:LEU:HD21	2:H:182:GLN:HE21	1.65	0.60
2:B:149:LEU:HD11	2:B:199:ARG:HG2	1.84	0.60
1:A:169:TRP:CD1	1:A:181:MET:HG3	2.36	0.60
1:C:169:TRP:CD1	1:C:181:MET:HG3	2.36	0.59
1:J:196:ASN:HD21	1:J:218:ASN:HB2	1.67	0.59
3:I:218:VAL:HG11	3:I:237:LEU:HD13	1.85	0.59
3:L:318:ALA:HB1	3:L:319:ILE:HG23	1.85	0.59
3:L:225:ILE:HD12	10:L:1098:HOH:O	2.02	0.58
1:J:169:TRP:CD1	1:J:181:MET:HG3	2.37	0.58
1:J:113:LYS:HA	1:J:146:TYR:OH	2.04	0.58
1:A:19:VAL:HG13	1:A:81:ILE:HB	1.84	0.58
1:A:21:MET:HG2	1:A:108:THR:HG21	1.84	0.57
2:K:199:ARG:CG	2:K:199:ARG:NH1	2.58	0.57
1:A:156:ILE:HD11	1:A:185:LEU:HD21	1.86	0.57
3:E:218:VAL:HG11	3:E:237:LEU:HD13	1.86	0.57
3:I:271:VAL:HG13	3:I:299:VAL:HG23	1.86	0.56
3:L:278:GLU:HG3	3:L:281:ARG:HH11	1.70	0.56
3:L:218:VAL:HG11	3:L:237:LEU:HD13	1.87	0.56
2:K:14:VAL:HG21	2:K:88:LEU:HD13	1.89	0.55
2:B:40:LYS:HE2	2:B:42:ARG:HD2	1.88	0.55
2:B:48:GLU:HA	4:B:607:EDO:H21	1.88	0.55
3:L:269:ILE:HG21	3:L:305:LEU:HD11	1.89	0.55
3:E:311:GLN:H	3:E:311:GLN:HE21	1.55	0.54
3:G:166:LYS:NZ	3:G:320:GLU:HG2	2.22	0.54
1:J:31:TYR:CZ	1:J:33:SER:HB2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:311:GLN:HE21	3:G:311:GLN:H	1.55	0.54
3:G:148:HIS:HD2	3:G:151:ARG:NH2	2.04	0.54
1:J:114:ARG:HG2	1:J:177:SER:HB2	1.90	0.54
2:B:214:SER:HA	5:B:704:GOL:H2	1.89	0.54
2:H:159:GLU:HG3	2:H:160:PRO:HA	1.90	0.54
2:B:159:GLU:HG3	2:B:160:PRO:HA	1.90	0.54
2:B:199:ARG:NH1	2:B:199:ARG:CG	2.55	0.53
2:D:174:VAL:HG22	2:D:192:VAL:HG23	1.90	0.53
2:K:159:GLU:HG3	2:K:160:PRO:HA	1.91	0.53
3:L:160:VAL:HG23	3:L:309:GLN:HE21	1.74	0.53
1:F:31:TYR:CZ	1:F:33:SER:HB2	2.43	0.53
2:D:199:ARG:CG	2:D:199:ARG:NH1	2.60	0.53
2:D:159:GLU:HG3	2:D:160:PRO:HA	1.92	0.52
2:H:220:LYS:HB3	2:K:214:SER:HB3	1.90	0.52
2:B:14:VAL:HG21	2:B:88:LEU:HD13	1.91	0.52
2:K:149:LEU:HD11	2:K:199:ARG:HG2	1.92	0.52
1:C:31:TYR:CZ	1:C:33:SER:HB2	2.45	0.52
3:L:311:GLN:H	3:L:311:GLN:HE21	1.56	0.52
3:G:240:ILE:HG12	3:G:269:ILE:HD12	1.92	0.52
3:I:311:GLN:HE21	3:I:311:GLN:H	1.55	0.52
3:I:170:LEU:HD13	3:I:189:PHE:HD2	1.74	0.52
1:F:19:VAL:HG13	1:F:81:ILE:HB	1.92	0.51
2:H:142:GLN:HE21	2:H:144:ASN:HD21	1.58	0.51
1:A:31:TYR:CZ	1:A:33:SER:HB2	2.44	0.51
2:D:14:VAL:HG21	2:D:88:LEU:HD13	1.91	0.51
1:C:17:GLU:O	1:C:84:VAL:HG23	2.11	0.51
2:H:14:VAL:HG21	2:H:88:LEU:HD13	1.93	0.51
2:D:104:TYR:HD1	3:E:151:ARG:HH21	1.58	0.51
2:K:34:ILE:HG22	2:K:102:GLY:HA2	1.94	0.50
3:L:209:THR:HG22	3:L:245:LYS:HA	1.93	0.49
2:K:6:LEU:HD23	2:K:26:PRO:HB3	1.94	0.49
2:D:199:ARG:HD2	2:D:200:PRO:HA	1.93	0.49
3:I:216:ARG:HB2	3:I:255:VAL:HG12	1.94	0.49
2:H:6:LEU:HD23	2:H:26:PRO:HB3	1.95	0.49
2:D:62:TYR:HD1	4:D:604:EDO:H12	1.77	0.49
2:B:182:GLN:O	2:B:183:SER:HB3	2.13	0.49
2:D:6:LEU:HD23	2:D:26:PRO:HB3	1.94	0.48
2:B:199:ARG:HD2	2:B:200:PRO:HA	1.96	0.48
2:D:76:THR:H	4:D:614:EDO:H22	1.78	0.48
3:I:166:LYS:HD2	3:I:317:PHE:HE1	1.79	0.48
2:K:7:GLN:HG3	5:K:707:GOL:H32	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:240:ILE:HG12	3:L:269:ILE:HD12	1.96	0.48
1:C:156:ILE:HD11	1:C:185:LEU:HD21	1.96	0.48
1:A:114:ARG:HD3	1:A:115:ALA:O	2.14	0.48
2:H:34:ILE:HD13	2:H:100:SER:HB2	1.95	0.48
2:B:6:LEU:HD23	2:B:26:PRO:HB3	1.95	0.47
2:B:195:PRO:HB2	2:B:198:PRO:CD	2.45	0.47
3:L:256:ILE:HG23	3:L:266:ARG:HH21	1.80	0.47
3:E:312:LEU:O	3:E:316:ILE:HG12	2.14	0.47
2:D:195:PRO:HB2	2:D:198:PRO:CD	2.45	0.47
3:G:278:GLU:O	3:G:282:GLN:HG3	2.15	0.46
2:D:82:TYR:CE1	2:H:202:GLU:HG2	2.50	0.46
2:H:142:GLN:HE21	2:H:144:ASN:ND2	2.14	0.45
1:J:97:TYR:HA	1:J:102:LEU:HD22	1.99	0.45
1:A:1:ASP:N	5:A:700:GOL:H2	2.31	0.45
3:I:312:LEU:O	3:I:316:ILE:HG12	2.17	0.45
3:G:172:SER:OG	3:G:222:LEU:HD22	2.16	0.45
2:H:195:PRO:HB2	2:H:198:PRO:CD	2.45	0.45
1:F:97:TYR:HA	1:F:102:LEU:HD22	1.99	0.45
1:C:97:TYR:HA	1:C:102:LEU:HD22	1.99	0.45
1:A:97:TYR:HA	1:A:102:LEU:HD22	1.98	0.44
2:B:49:TRP:H	4:B:607:EDO:H21	1.82	0.44
2:K:195:PRO:HB2	2:K:198:PRO:CD	2.46	0.44
1:C:69:THR:HG22	10:C:1239:HOH:O	2.18	0.44
2:D:61:LYS:HA	4:D:604:EDO:H21	2.00	0.44
3:L:256:ILE:HG23	3:L:266:ARG:NH2	2.33	0.44
3:L:133:SER:O	3:L:169:THR:HA	2.18	0.44
2:H:130:PRO:HB3	2:H:156:TYR:HB3	2.00	0.43
3:E:160:VAL:HG23	3:E:309:GLN:HE21	1.83	0.43
3:I:132:ASP:HA	3:I:168:LYS:HB3	2.00	0.43
1:F:181:MET:HE3	1:F:183:SER:HB2	2.00	0.43
1:F:95:GLN:HE21	1:F:95:GLN:HB3	1.67	0.43
3:G:217:LYS:NZ	5:G:703:GOL:O2	2.50	0.43
2:K:93:THR:HG23	2:K:121:THR:HA	2.00	0.43
2:B:93:THR:HG23	2:B:121:THR:HA	2.01	0.43
2:D:42:ARG:HG3	2:D:43:PRO:HD2	2.01	0.43
2:H:121:THR:HG21	2:H:186:TYR:OH	2.19	0.43
1:J:129:GLU:HG2	2:K:219:LYS:HZ1	1.84	0.42
2:B:49:TRP:H	4:B:607:EDO:C2	2.32	0.42
3:G:132:ASP:HA	3:G:168:LYS:HB3	2.00	0.42
3:L:132:ASP:HA	3:L:168:LYS:HB3	2.01	0.42
2:K:34:ILE:HD13	2:K:100:SER:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:153:MET:O	3:G:157:VAL:HG23	2.20	0.42
1:J:67:ARG:HG3	1:J:81:ILE:HG23	2.01	0.42
2:H:93:THR:HG23	2:H:121:THR:HA	2.01	0.42
2:H:175:HIS:HB2	2:H:191:SER:HB3	2.00	0.42
2:B:136:ALA:HB3	2:B:224:ARG:HH21	1.85	0.42
1:A:2:ILE:HG12	1:A:27:GLN:HB2	2.02	0.42
2:B:82:TYR:CE1	2:K:202:GLU:HG2	2.54	0.42
3:G:272:GLY:HA2	3:G:300:ASN:O	2.20	0.42
3:L:219:VAL:HG22	3:L:264:VAL:HG21	2.01	0.41
1:F:196:ASN:ND2	1:F:218:ASN:H	2.18	0.41
3:G:138:LEU:HB2	3:G:218:VAL:HG21	2.01	0.41
1:J:181:MET:HE3	1:J:183:SER:HB2	2.02	0.41
2:D:93:THR:HG23	2:D:121:THR:HA	2.02	0.41
1:A:196:ASN:ND2	1:A:218:ASN:H	2.18	0.41
2:D:130:PRO:HB3	2:D:156:TYR:HB3	2.01	0.41
1:C:67:ARG:HG3	1:C:81:ILE:HG23	2.03	0.41
3:I:140:ASP:HB2	3:I:241:THR:HA	2.02	0.41
3:I:311:GLN:H	3:I:311:GLN:NE2	2.18	0.41
2:K:174:VAL:HG22	2:K:192:VAL:HG23	2.02	0.41
1:J:2:ILE:HG12	1:J:27:GLN:HB2	2.03	0.41
1:F:67:ARG:HB3	10:F:3223:HOH:O	2.21	0.41
3:E:132:ASP:HA	3:E:168:LYS:HB3	2.02	0.41
1:C:220:CYS:HB3	2:D:226:CYS:HB3	1.75	0.41
1:F:114:ARG:HD3	1:F:115:ALA:O	2.20	0.41
3:G:141:GLY:HA3	3:G:175:GLN:OE1	2.21	0.41
3:G:200:LYS:HE3	2:B:28:GLY:HA2	2.02	0.41
3:L:311:GLN:NE2	3:L:311:GLN:H	2.19	0.40
1:A:181:MET:HE3	1:A:183:SER:HB2	2.04	0.40
1:C:67:ARG:HB2	1:C:82:SER:O	2.21	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	218/220 (99%)	212 (97%)	4 (2%)	2 (1%)	21	49
1	C	218/220 (99%)	212 (97%)	5 (2%)	1 (0%)	34	63
1	F	218/220 (99%)	211 (97%)	7 (3%)	0	100	100
1	J	218/220 (99%)	211 (97%)	6 (3%)	1 (0%)	34	63
2	B	222/224 (99%)	210 (95%)	10 (4%)	2 (1%)	21	49
2	D	222/224 (99%)	212 (96%)	6 (3%)	4 (2%)	11	27
2	H	222/224 (99%)	211 (95%)	10 (4%)	1 (0%)	34	63
2	K	222/224 (99%)	213 (96%)	8 (4%)	1 (0%)	34	63
3	E	187/190 (98%)	178 (95%)	7 (4%)	2 (1%)	17	42
3	G	187/190 (98%)	179 (96%)	7 (4%)	1 (0%)	34	63
3	I	187/190 (98%)	180 (96%)	5 (3%)	2 (1%)	17	42
3	L	187/190 (98%)	178 (95%)	8 (4%)	1 (0%)	34	63
All	All	2508/2536 (99%)	2407 (96%)	83 (3%)	18 (1%)	26	55

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	225	ASP
3	E	168	LYS
2	H	141	ALA
1	J	46	PRO
2	B	142	GLN
2	D	142	GLN
3	G	177	SER
1	A	219	GLU
2	B	183	SER
3	E	177	SER
3	I	177	SER
3	L	177	SER
1	C	83	SER
2	D	183	SER
1	A	83	SER
2	K	182	GLN
3	I	319	ILE
2	D	143	THR

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	197/197 (100%)	180 (91%)	17 (9%)	13	29
1	C	197/197 (100%)	183 (93%)	14 (7%)	18	41
1	F	197/197 (100%)	179 (91%)	18 (9%)	12	26
1	J	197/197 (100%)	175 (89%)	22 (11%)	7	17
2	B	192/192 (100%)	181 (94%)	11 (6%)	25	53
2	D	192/192 (100%)	181 (94%)	11 (6%)	25	53
2	H	192/192 (100%)	184 (96%)	8 (4%)	36	68
2	K	192/192 (100%)	183 (95%)	9 (5%)	32	63
3	E	169/169 (100%)	156 (92%)	13 (8%)	16	36
3	G	169/169 (100%)	163 (96%)	6 (4%)	42	73
3	I	169/169 (100%)	159 (94%)	10 (6%)	24	51
3	L	169/169 (100%)	160 (95%)	9 (5%)	28	57
All	All	2232/2232 (100%)	2084 (93%)	148 (7%)	21	45

All (148) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	5	THR
1	C	13	VAL
1	C	14	SER
1	C	39	LEU
1	C	54	ILE
1	C	78	THR
1	C	95	GLN
1	C	111	GLU
1	C	114	ARG
1	C	151	ASN
1	C	152	VAL
1	C	160	GLU
1	C	199	THR
1	C	203	THR

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Mol	Chain	Res	Type
2	D	26	PRO
2	D	30	ASN
2	D	42	ARG
2	D	52	ARG
2	D	139	SER
2	D	161	VAL
2	D	171	SER
2	D	196	SER
2	D	197	SER
2	D	199	ARG
2	D	207	ASN
3	G	155	GLU
3	G	225	ILE
3	G	226	THR
3	G	230	ARG
3	G	242	ASP
3	G	311	GLN
1	A	5	THR
1	A	13	VAL
1	A	14	SER
1	A	19	VAL
1	A	26	SER
1	A	27	GLN
1	A	39	LEU
1	A	54	ILE
1	A	78	THR
1	A	84	VAL
1	A	95	GLN
1	A	111	GLU
1	A	114	ARG
1	A	151	ASN
1	A	152	VAL
1	A	199	THR
1	A	203	THR
2	B	13	LEU
2	B	30	ASN
2	B	42	ARG
2	B	52	ARG
2	B	161	VAL
2	B	184	ASP
2	B	196	SER
2	B	197	SER

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Mol	Chain	Res	Type
2	B	199	ARG
2	B	207	ASN
2	B	224	ARG
3	E	155	GLU
3	E	158	SER
3	E	172	SER
3	E	225	ILE
3	E	226	THR
3	E	230	ARG
3	E	242	ASP
3	E	245	LYS
3	E	290	LYS
3	E	309	GLN
3	E	311	GLN
3	E	319	ILE
3	E	320	GLU
1	F	5	THR
1	F	13	VAL
1	F	14	SER
1	F	19	VAL
1	F	27	GLN
1	F	39	LEU
1	F	54	ILE
1	F	78	THR
1	F	84	VAL
1	F	95	GLN
1	F	111	GLU
1	F	114	ARG
1	F	151	ASN
1	F	152	VAL
1	F	160	GLU
1	F	199	THR
1	F	203	THR
1	F	220	CYS
2	H	30	ASN
2	H	44	GLU
2	H	52	ARG
2	H	161	VAL
2	H	196	SER
2	H	197	SER
2	H	199	ARG
2	H	207	ASN

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Mol	Chain	Res	Type
3	I	155	GLU
3	I	158	SER
3	I	172	SER
3	I	225	ILE
3	I	226	THR
3	I	230	ARG
3	I	242	ASP
3	I	279	LYS
3	I	300	ASN
3	I	311	GLN
1	J	3	GLU
1	J	5	THR
1	J	13	VAL
1	J	14	SER
1	J	19	VAL
1	J	26	SER
1	J	27	GLN
1	J	39	LEU
1	J	46	PRO
1	J	54	ILE
1	J	78	THR
1	J	84	VAL
1	J	95	GLN
1	J	111	GLU
1	J	114	ARG
1	J	151	ASN
1	J	152	VAL
1	J	160	GLU
1	J	196	ASN
1	J	199	THR
1	J	203	THR
1	J	220	CYS
2	K	30	ASN
2	K	52	ARG
2	K	143	THR
2	K	161	VAL
2	K	184	ASP
2	K	196	SER
2	K	197	SER
2	K	199	ARG
2	K	226	CYS
3	L	158	SER

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Mol	Chain	Res	Type
3	L	225	ILE
3	L	226	THR
3	L	230	ARG
3	L	242	ASP
3	L	273	ASP
3	L	309	GLN
3	L	311	GLN
3	L	319	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	35	GLN
1	C	44	GLN
1	C	95	GLN
1	C	167	ASN
2	D	41	GLN
3	G	148	HIS
3	G	224	ASN
3	G	227	ASN
3	G	282	GLN
3	G	311	GLN
1	A	27	GLN
1	A	35	GLN
1	A	44	GLN
1	A	95	GLN
1	A	167	ASN
1	A	196	ASN
1	A	216	ASN
2	B	41	GLN
2	B	103	HIS
3	E	224	ASN
3	E	309	GLN
3	E	311	GLN
1	F	27	GLN
1	F	35	GLN
1	F	44	GLN
1	F	95	GLN
1	F	167	ASN
2	H	41	GLN
2	H	144	ASN
2	H	175	HIS

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Mol	Chain	Res	Type
2	H	182	GLN
3	I	210	HIS
3	I	224	ASN
3	I	227	ASN
3	I	309	GLN
3	I	311	GLN
1	J	27	GLN
1	J	35	GLN
1	J	44	GLN
1	J	95	GLN
1	J	167	ASN
1	J	196	ASN
2	K	41	GLN
3	L	224	ASN
3	L	282	GLN
3	L	309	GLN
3	L	311	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 49 ligands modelled in this entry, 12 are monoatomic - leaving 37 for Mogul analysis.

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
4	EDO	A	618	-	3,3,3	0.50	0	2,2,2	0.43	0
4	EDO	A	625	-	3,3,3	0.46	0	2,2,2	0.39	0
5	GOL	A	700	-	5,5,5	0.73	0	5,5,5	0.77	0
4	EDO	B	606	-	3,3,3	0.50	0	2,2,2	0.47	0
4	EDO	B	607	-	3,3,3	0.62	0	2,2,2	0.10	0
4	EDO	B	616	-	3,3,3	0.54	0	2,2,2	0.18	0
4	EDO	B	624	-	3,3,3	0.51	0	2,2,2	0.26	0
5	GOL	B	701	-	5,5,5	0.59	0	5,5,5	0.61	0
5	GOL	B	704	-	5,5,5	0.72	0	5,5,5	0.45	0
4	EDO	C	609	-	3,3,3	0.59	0	2,2,2	0.19	0
4	EDO	C	628	-	3,3,3	0.49	0	2,2,2	0.56	0
5	GOL	C	710	-	5,5,5	1.26	0	5,5,5	1.35	1 (20%)
4	EDO	D	604	-	3,3,3	0.65	0	2,2,2	0.06	0
4	EDO	D	605	-	3,3,3	0.54	0	2,2,2	0.28	0
4	EDO	D	613	-	3,3,3	0.56	0	2,2,2	0.21	0
4	EDO	D	614	-	3,3,3	0.46	0	2,2,2	0.61	0
4	EDO	D	619	-	3,3,3	0.51	0	2,2,2	0.42	0
5	GOL	D	702	-	5,5,5	0.69	0	5,5,5	0.97	1 (20%)
5	GOL	D	705	-	5,5,5	1.01	0	5,5,5	0.73	0
5	GOL	D	708	-	5,5,5	1.14	0	5,5,5	0.80	0
4	EDO	E	600	-	3,3,3	0.53	0	2,2,2	0.31	0
4	EDO	E	627	-	3,3,3	0.51	0	2,2,2	0.38	0
4	EDO	F	621	-	3,3,3	0.58	0	2,2,2	0.27	0
4	EDO	G	620	-	3,3,3	0.45	0	2,2,2	0.47	0
5	GOL	G	703	-	5,5,5	0.96	0	5,5,5	1.18	0
9	PEG	H	1000	-	6,6,6	0.46	0	5,5,5	0.51	0
4	EDO	J	603	-	3,3,3	0.65	0	2,2,2	0.10	0
4	EDO	J	622	-	3,3,3	0.58	0	2,2,2	0.11	0
5	GOL	J	709	-	5,5,5	1.00	0	5,5,5	0.92	0
4	EDO	K	601	-	3,3,3	0.52	0	2,2,2	0.43	0
4	EDO	K	608	-	3,3,3	0.52	0	2,2,2	0.36	0
4	EDO	K	611	-	3,3,3	0.56	0	2,2,2	0.25	0
4	EDO	K	612	-	3,3,3	0.61	0	2,2,2	0.31	0
5	GOL	K	706	-	5,5,5	0.69	0	5,5,5	0.87	0
5	GOL	K	707	-	5,5,5	0.55	0	5,5,5	0.84	0
4	EDO	L	610	-	3,3,3	0.54	0	2,2,2	0.36	0
4	EDO	L	615	-	3,3,3	0.47	0	2,2,2	0.45	0

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
4	EDO	A	618	-	-	0/1/1/1	0/0/0/0
4	EDO	A	625	-	-	0/1/1/1	0/0/0/0
5	GOL	A	700	-	-	0/4/4/4	0/0/0/0
4	EDO	B	606	-	-	0/1/1/1	0/0/0/0
4	EDO	B	607	-	-	0/1/1/1	0/0/0/0
4	EDO	B	616	-	-	0/1/1/1	0/0/0/0
4	EDO	B	624	-	-	0/1/1/1	0/0/0/0
5	GOL	B	701	-	-	0/4/4/4	0/0/0/0
5	GOL	B	704	-	-	0/4/4/4	0/0/0/0
4	EDO	C	609	-	-	0/1/1/1	0/0/0/0
4	EDO	C	628	-	-	0/1/1/1	0/0/0/0
5	GOL	C	710	-	-	0/4/4/4	0/0/0/0
4	EDO	D	604	-	-	0/1/1/1	0/0/0/0
4	EDO	D	605	-	-	0/1/1/1	0/0/0/0
4	EDO	D	613	-	-	0/1/1/1	0/0/0/0
4	EDO	D	614	-	-	0/1/1/1	0/0/0/0
4	EDO	D	619	-	-	0/1/1/1	0/0/0/0
5	GOL	D	702	-	-	0/4/4/4	0/0/0/0
5	GOL	D	705	-	-	0/4/4/4	0/0/0/0
5	GOL	D	708	-	-	0/4/4/4	0/0/0/0
4	EDO	E	600	-	-	0/1/1/1	0/0/0/0
4	EDO	E	627	-	-	0/1/1/1	0/0/0/0
4	EDO	F	621	-	-	0/1/1/1	0/0/0/0
4	EDO	G	620	-	-	0/1/1/1	0/0/0/0
5	GOL	G	703	-	-	0/4/4/4	0/0/0/0
9	PEG	H	1000	-	-	0/4/4/4	0/0/0/0
4	EDO	J	603	-	-	0/1/1/1	0/0/0/0
4	EDO	J	622	-	-	0/1/1/1	0/0/0/0
5	GOL	J	709	-	-	0/4/4/4	0/0/0/0
4	EDO	K	601	-	-	0/1/1/1	0/0/0/0
4	EDO	K	608	-	-	0/1/1/1	0/0/0/0
4	EDO	K	611	-	-	0/1/1/1	0/0/0/0
4	EDO	K	612	-	-	0/1/1/1	0/0/0/0
5	GOL	K	706	-	-	0/4/4/4	0/0/0/0
5	GOL	K	707	-	-	0/4/4/4	0/0/0/0
4	EDO	L	610	-	-	0/1/1/1	0/0/0/0
4	EDO	L	615	-	-	0/1/1/1	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	702	GOL	O1-C1-C2	2.09	120.30	110.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	710	GOL	O2-C2-C3	2.63	120.69	108.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	700	GOL	1	0
4	B	607	EDO	3	0
5	B	704	GOL	1	0
4	D	604	EDO	2	0
4	D	614	EDO	1	0
5	D	702	GOL	1	0
4	E	600	EDO	1	0
5	G	703	GOL	1	0
5	K	707	GOL	1	0

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 ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	220/220 (100%)	0.07	2 (0%) 85 86	25, 42, 66, 94	0
1	C	220/220 (100%)	0.09	3 (1%) 78 77	26, 41, 80, 113	0
1	F	220/220 (100%)	1.05	35 (15%) 3 2	38, 74, 97, 112	0
1	J	220/220 (100%)	0.11	2 (0%) 85 86	26, 46, 66, 110	0
2	B	224/224 (100%)	0.18	10 (4%) 37 36	24, 39, 73, 105	0
2	D	224/224 (100%)	-0.01	3 (1%) 79 79	21, 37, 70, 113	0
2	H	224/224 (100%)	0.12	6 (2%) 58 58	28, 47, 77, 111	0
2	K	224/224 (100%)	0.15	7 (3%) 52 52	24, 41, 74, 111	0
3	E	189/190 (99%)	0.23	6 (3%) 51 51	28, 53, 76, 115	0
3	G	189/190 (99%)	0.12	4 (2%) 67 68	29, 44, 73, 104	0
3	I	189/190 (99%)	1.41	55 (29%) 1 0	42, 87, 116, 125	0
3	L	189/190 (99%)	1.09	35 (18%) 2 1	30, 73, 112, 122	0
All	All	2532/2536 (99%)	0.37	168 (6%) 22 20	21, 48, 98, 125	0

All (168) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	L	167	SER	6.4
3	L	297	PHE	6.4
3	L	319	ILE	5.9
2	K	3	GLN	5.6
2	B	3	GLN	5.4
1	A	220	CYS	5.4
2	D	3	GLN	5.3
3	I	167	SER	5.3
1	J	220	CYS	5.3
2	K	226	CYS	5.3
3	E	320	GLU	5.1

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Mol	Chain	Res	Type	RSRZ
2	K	4	VAL	5.0
3	L	316	ILE	5.0
3	I	259	ALA	4.9
3	I	225	ILE	4.9
1	F	205	LYS	4.8
3	I	172	SER	4.6
3	I	222	LEU	4.6
3	I	219	VAL	4.5
3	I	236	ILE	4.5
1	C	220	CYS	4.4
3	I	227	ASN	4.3
3	E	319	ILE	4.2
3	I	316	ILE	4.2
2	B	225	ASP	4.2
2	D	226	CYS	4.2
3	L	314	GLU	4.1
3	L	267	TYR	4.1
1	F	200	CYS	4.1
3	I	318	ALA	4.0
3	I	234	PHE	3.9
3	I	294	ASP	3.8
3	I	134	ASP	3.7
1	F	212	VAL	3.6
3	L	236	ILE	3.6
1	F	211	ILE	3.6
3	I	312	LEU	3.5
3	L	320	GLU	3.5
3	L	318	ALA	3.5
1	F	206	THR	3.4
1	F	84	VAL	3.4
1	F	112	LEU	3.4
3	I	293	ARG	3.4
3	I	256	ILE	3.4
3	I	237	LEU	3.3
3	L	165	LYS	3.3
3	I	228	GLY	3.3
2	H	3	GLN	3.3
3	I	168	LYS	3.3
3	L	168	LYS	3.3
3	L	281	ARG	3.2
1	F	218	ASN	3.2
2	B	226	CYS	3.2

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Mol	Chain	Res	Type	RSRZ
3	I	288	ALA	3.2
3	I	223	PHE	3.2
3	L	134	ASP	3.1
3	I	164	LEU	3.1
3	I	267	TYR	3.1
3	L	137	PHE	3.1
2	B	224	ARG	3.0
2	H	140	ALA	3.0
3	I	133	SER	3.0
3	L	307	THR	3.0
3	I	232	ASN	3.0
3	I	319	ILE	3.0
1	F	146	TYR	2.9
3	I	276	ARG	2.9
1	F	88	ASP	2.9
1	F	81	ILE	2.9
3	L	164	LEU	2.9
1	F	159	SER	2.9
1	F	86	ALA	2.9
3	G	320	GLU	2.9
1	F	208	THR	2.9
3	I	218	VAL	2.8
2	B	141	ALA	2.8
3	I	235	LYS	2.8
3	I	292	PRO	2.8
3	I	166	LYS	2.8
3	I	169	THR	2.8
2	H	43	PRO	2.7
1	F	82	SER	2.7
3	L	312	LEU	2.7
2	H	226	CYS	2.7
3	L	313	ARG	2.7
2	K	142	GLN	2.7
1	F	38	TYR	2.7
3	L	287	ILE	2.7
2	B	4	VAL	2.7
3	I	314	GLU	2.7
3	I	261	ARG	2.6
1	F	19	VAL	2.6
2	H	4	VAL	2.6
3	I	132	ASP	2.6
1	F	13	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
3	I	287	ILE	2.6
1	F	158	GLY	2.6
1	F	3	GLU	2.6
1	F	220	CYS	2.6
3	I	226	THR	2.6
1	F	155	LYS	2.6
3	L	317	PHE	2.6
3	L	278	GLU	2.6
3	I	265	ILE	2.5
3	L	308	ILE	2.5
1	C	187	LEU	2.5
2	K	144	ASN	2.5
3	I	264	VAL	2.5
2	K	225	ASP	2.5
3	L	280	SER	2.5
3	I	171	PHE	2.5
3	I	257	PRO	2.5
1	F	23	CYS	2.5
3	L	295	HIS	2.5
1	F	187	LEU	2.5
2	B	140	ALA	2.4
3	I	255	VAL	2.4
3	I	268	VAL	2.4
1	C	194	ARG	2.4
1	F	111	GLU	2.4
3	L	268	VAL	2.4
3	I	216	ARG	2.4
3	L	135	ILE	2.4
3	I	295	HIS	2.4
3	G	167	SER	2.4
1	F	154	TRP	2.4
1	J	46	PRO	2.4
3	I	311	GLN	2.4
3	I	290	LYS	2.4
3	E	234	PHE	2.4
3	L	265	ILE	2.4
3	L	294	ASP	2.4
2	H	142	GLN	2.4
3	L	194	ASN	2.4
3	L	298	GLN	2.3
1	F	10	SER	2.3
1	F	121	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	79	LEU	2.3
3	L	230	ARG	2.3
1	F	115	ALA	2.2
3	L	255	VAL	2.2
1	F	8	PRO	2.2
3	L	218	VAL	2.2
1	F	18	LYS	2.2
2	K	182	GLN	2.2
3	I	281	ARG	2.2
1	A	218	ASN	2.2
3	G	232	ASN	2.2
3	I	220	ARG	2.2
1	F	162	GLN	2.2
3	E	164	LEU	2.2
2	B	142	GLN	2.2
3	I	317	PHE	2.2
3	I	238	VAL	2.1
3	I	135	ILE	2.1
3	G	231	LYS	2.1
3	E	231	LYS	2.1
3	L	166	LYS	2.1
3	E	167	SER	2.1
1	F	147	PRO	2.1
3	I	269	ILE	2.0
2	B	44	GLU	2.0
2	B	139	SER	2.0
3	I	231	LYS	2.0
3	I	250	LEU	2.0
3	L	234	PHE	2.0
1	F	179	TYR	2.0
2	D	143	THR	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	GOL	B	704	6/6	0.82	0.36	8.85	54,56,56,57	0
4	EDO	D	614	4/4	0.85	0.26	6.80	49,49,50,51	0
4	EDO	C	628	4/4	0.91	0.21	5.49	50,50,52,56	0
5	GOL	K	707	6/6	0.83	0.34	4.85	66,68,68,69	0
4	EDO	A	618	4/4	0.72	0.32	4.73	57,59,62,62	0
4	EDO	B	616	4/4	0.77	0.25	4.20	63,64,64,65	0
4	EDO	E	600	4/4	0.88	0.24	4.20	42,43,48,49	0
4	EDO	B	607	4/4	0.84	0.28	3.81	47,50,51,55	0
4	EDO	D	613	4/4	0.87	0.23	3.46	51,51,55,58	0
4	EDO	C	609	4/4	0.81	0.24	3.22	54,54,56,57	0
5	GOL	J	709	6/6	0.79	0.30	2.53	59,60,61,62	0
4	EDO	D	619	4/4	0.85	0.22	2.53	53,55,57,60	0
5	GOL	K	706	6/6	0.84	0.25	2.29	60,62,62,62	0
5	GOL	A	700	6/6	0.69	0.28	2.15	68,69,70,71	0
4	EDO	E	627	4/4	0.91	0.20	2.04	54,57,60,62	0
4	EDO	G	620	4/4	0.96	0.21	2.02	38,38,39,46	0
5	GOL	G	703	6/6	0.93	0.20	1.45	48,48,49,49	0
4	EDO	D	605	4/4	0.87	0.22	1.29	49,50,50,55	0
4	EDO	K	612	4/4	0.87	0.21	1.23	39,40,41,42	0
9	PEG	H	1000	7/7	0.86	0.26	1.21	59,59,62,63	0
5	GOL	D	705	6/6	0.83	0.20	1.10	59,61,61,62	0
5	GOL	D	702	6/6	0.75	0.18	1.06	61,63,64,64	0
4	EDO	J	603	4/4	0.91	0.18	0.85	40,42,45,45	0
4	EDO	F	621	4/4	0.86	0.22	0.83	56,57,58,62	0
5	GOL	D	708	6/6	0.83	0.19	0.40	50,50,51,51	0
4	EDO	K	601	4/4	0.90	0.17	0.24	40,40,41,46	0
5	GOL	C	710	6/6	0.84	0.19	0.10	53,56,57,58	0
4	EDO	L	615	4/4	0.88	0.24	0.06	51,54,55,55	0
4	EDO	K	611	4/4	0.78	0.20	-0.18	60,62,62,65	0
8	CL	J	903	1/1	0.72	0.18	-0.21	70,70,70,70	0
4	EDO	L	610	4/4	0.77	0.20	-0.34	57,58,58,65	0
4	EDO	B	606	4/4	0.89	0.15	-0.43	43,43,47,48	0
7	CA	E	500	1/1	0.96	0.17	-0.55	40,40,40,40	0
4	EDO	A	625	4/4	0.96	0.16	-0.78	41,43,45,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
7	CA	G	500	1/1	0.97	0.15	-1.86	27,27,27,27	0
7	CA	I	500	1/1	0.89	0.11	-2.38	57,57,57,57	0
7	CA	L	500	1/1	0.96	0.11	-3.22	40,40,40,40	0
4	EDO	D	604	4/4	0.84	0.26	-	37,40,41,48	0
6	NA	H	802	1/1	0.88	0.17	-	58,58,58,58	0
8	CL	K	904	1/1	0.79	0.10	-	63,63,63,63	0
5	GOL	B	701	6/6	0.84	0.36	-	71,71,72,73	0
6	NA	H	801	1/1	0.91	0.23	-	37,37,37,37	0
4	EDO	B	624	4/4	0.89	0.20	-	55,55,56,64	0
6	NA	D	800	1/1	0.93	0.17	-	41,41,41,41	0
8	CL	E	906	1/1	0.93	0.14	-	71,71,71,71	0
4	EDO	K	608	4/4	0.82	0.19	-	52,56,56,60	0
8	CL	G	907	1/1	0.90	0.13	-	74,74,74,74	0
4	EDO	J	622	4/4	0.86	0.24	-	55,56,57,66	0
8	CL	E	902	1/1	0.83	0.14	-	60,60,60,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.