



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 08:32 AM UTC

PDB ID : 4EFP / pdb_00004efp
Title : Bombyx mori lipoprotein 7 isolated from its natural source at 1.33 Å resolution
Authors : Pietrzyk, A.J.; Panjikar, S.; Bujacz, A.; Mueller-Dieckmann, J.; Jaskolski, M.;
Bujacz, G.
Deposited on : 2012-03-30
Resolution : 1.33 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)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

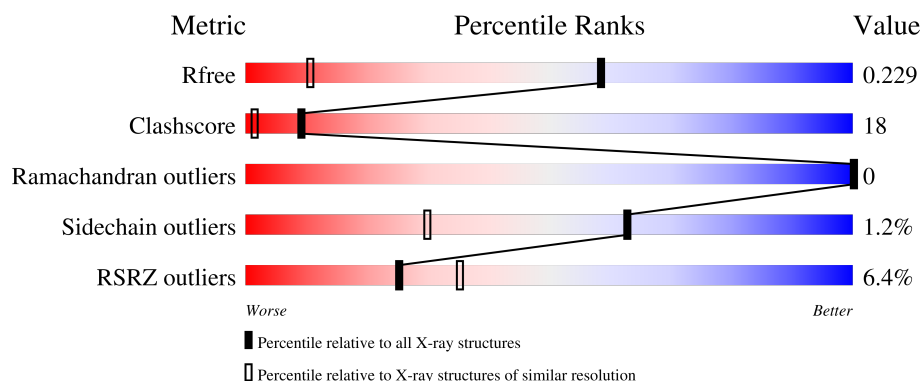
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.33 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}	180053	2194 (1.36-1.32)
Clashscore	190562	2222 (1.36-1.32)
Ramachandran outliers	187476	2197 (1.36-1.32)
Sidechain outliers	187428	2197 (1.36-1.32)
RSRZ outliers	180081	2193 (1.36-1.32)

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 $\leq 5\%$. 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	239	7% 77% 20% ..
1	B	239	5% 79% 17% ..

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
5	EDO	A	305	-	-	X	-
5	EDO	A	306	-	-	X	-

2 Entry composition [i](#)

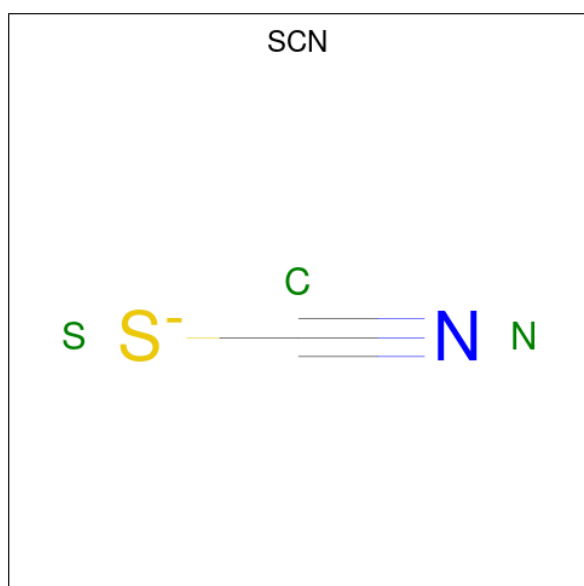
There are 8 unique types of molecules in this entry. The entry contains 4600 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 30kDa protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	S	0	16	0
			2010	1294	341	366	9			
1	B	235	Total	C	N	O	S	0	15	0
			1996	1289	336	361	10			

- Molecule 2 is THIOCYANATE ION (CCD ID: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	S	0	0
			3	1	1	1		
2	B	1	Total	C	N	S	0	0
			3	1	1	1		

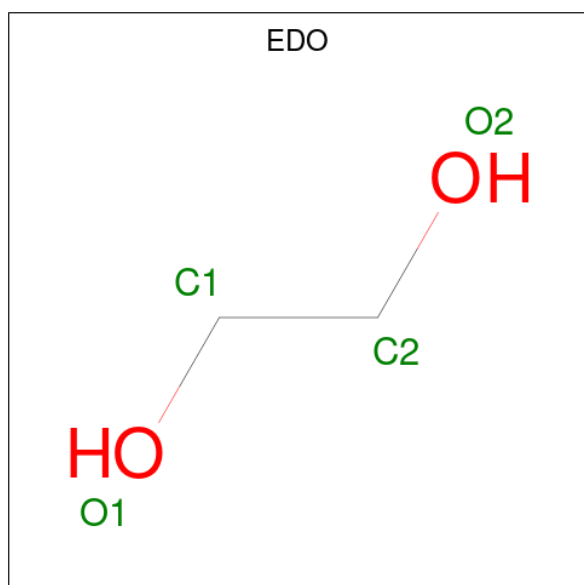
- Molecule 3 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cd 1 1	0	0
3	B	1	Total Cd 1 1	0	0

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

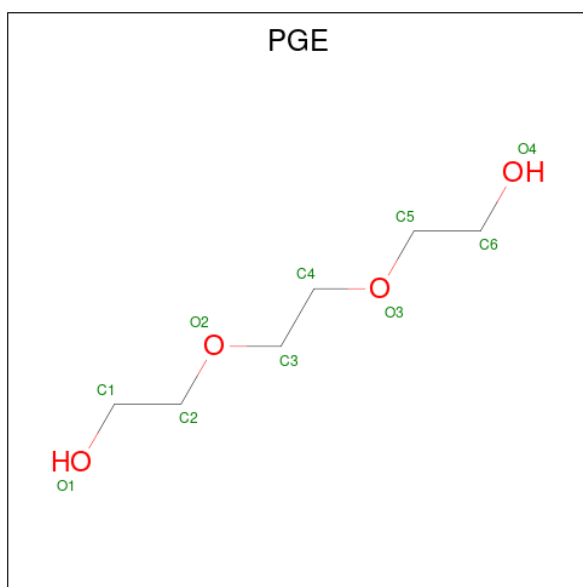
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total K 1 1	0	0
4	B	1	Total K 1 1	0	0

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



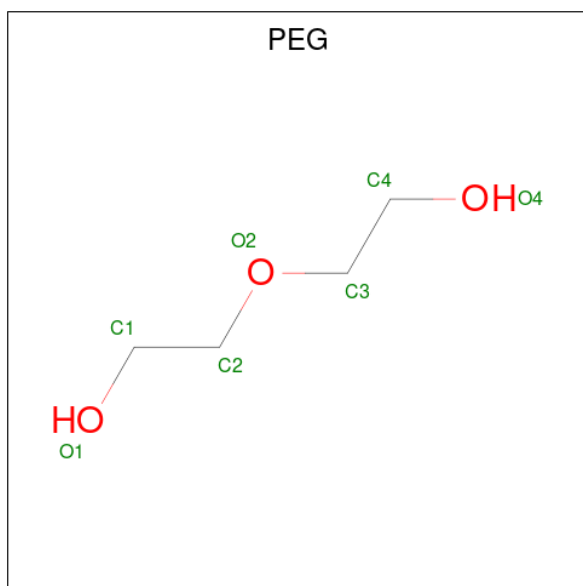
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	B	1	Total C O 4 2 2	0	0

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

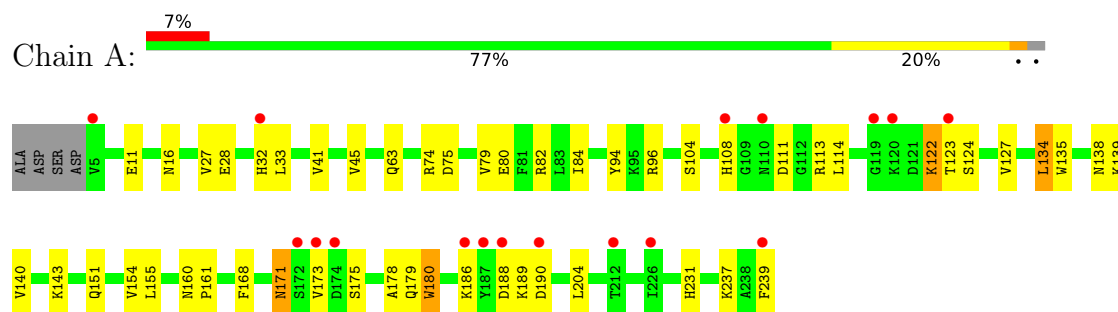
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	298	Total 298	O 298	0	0
8	B	253	Total 253	O 253	0	0

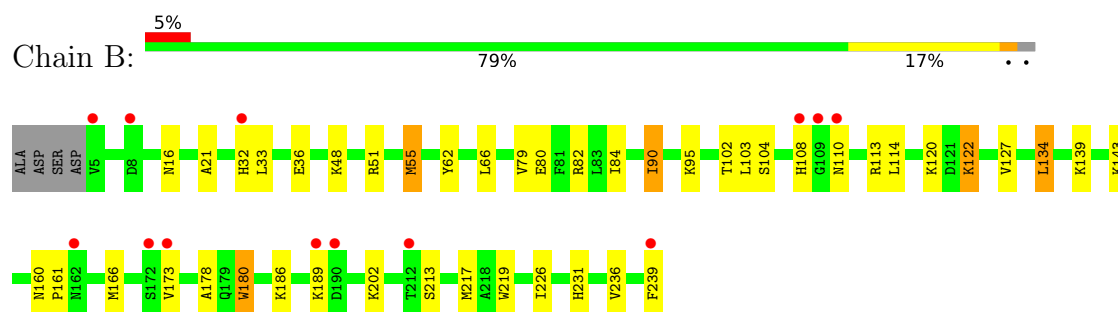
3 Residue-property plots

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 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: 30kDa protein



• Molecule 1: 30kDa protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	42.21Å 49.99Å 55.24Å 93.41° 94.70° 102.78°	Depositor
Resolution (Å)	23.00 – 1.33 23.00 – 1.33	Depositor EDS
% Data completeness (in resolution range)	93.6 (23.00-1.33) 94.0 (23.00-1.33)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 1.33Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.184 , 0.231 0.187 , 0.229	Depositor DCC
R_{free} test set	1223 reflections (1.21%)	wwPDB-VP
Wilson B-factor (Å ²)	11.6	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4600	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.51% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, 0AF, EDO, K, SCN, CD, PGE

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	1.01	2/2084 (0.1%)	1.01	4/2811 (0.1%)
1	B	1.05	1/2069 (0.0%)	1.04	0/2789
All	All	1.03	3/4153 (0.1%)	1.03	4/5600 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	140	VAL	CA-CB	8.01	1.63	1.54
1	B	90	ILE	C-O	5.43	1.29	1.23
1	A	94	TYR	C-O	-5.33	1.17	1.24

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	171	ASN	N-CA-C	9.23	122.53	111.82
1	A	122	LYS	N-CA-C	-7.12	102.91	112.94
1	A	188	ASP	CB-CA-C	-5.43	110.29	116.54
1	A	123	THR	N-CA-C	5.43	121.59	114.75

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	2010	0	2011	82	0
1	B	1996	0	2017	65	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	12	0	18	14	0
5	B	4	0	6	0	0
6	A	10	0	14	0	0
7	B	7	0	10	1	0
8	A	298	0	0	33	0
8	B	253	0	0	28	0
All	All	4600	0	4076	148	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 18.

All (148) 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:166[B]:MET:HG2	8:B:533:HOH:O	1.52	1.07
1:B:122:LYS:O	1:B:122:LYS:HG2	1.55	1.06
1:A:179:GLN:HE22	5:A:306:EDO:H22	1.19	1.04
1:A:179:GLN:NE2	5:A:306:EDO:H22	1.76	1.00
1:B:108:HIS:HB3	8:B:647:HOH:O	1.59	0.99
1:B:16:ASN:HB2	8:B:617:HOH:O	1.63	0.98
1:B:143[B]:LYS:HE3	1:B:173:VAL:HG11	1.44	0.97
1:B:189:LYS:HG3	8:B:607:HOH:O	1.65	0.97
1:B:62:TYR:CE2	1:B:66:LEU:HD11	2.01	0.96
1:A:108:HIS:HB3	8:A:627:HOH:O	1.67	0.95
1:A:63:GLN:HB3	8:A:656:HOH:O	1.69	0.93
1:B:80:GLU:O	1:B:84[B]:ILE:HD13	1.69	0.93
1:A:143:LYS:HE3	1:A:173:VAL:HG11	1.52	0.90
1:A:179:GLN:HE22	5:A:306:EDO:C2	1.85	0.89
1:A:139:LYS:HE2	8:A:611:HOH:O	1.71	0.89
1:A:96[B]:ARG:HD3	8:A:551:HOH:O	1.73	0.88
1:A:33:LEU:HG	8:A:679:HOH:O	1.72	0.88
1:B:62:TYR:CZ	1:B:66:LEU:HD11	2.11	0.86
1:B:21:ALA:CB	8:B:609:HOH:O	2.23	0.85
1:B:108:HIS:CE1	8:B:646:HOH:O	2.29	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:VAL:HG12	8:A:642:HOH:O	1.78	0.84
1:B:102[B]:THR:HG21	8:B:613:HOH:O	1.78	0.84
1:B:213:SER:HA	8:B:650:HOH:O	1.77	0.83
1:B:21:ALA:HB1	8:B:609:HOH:O	1.78	0.83
1:A:27:VAL:HG23	8:A:656:HOH:O	1.77	0.82
1:B:122:LYS:HG2	8:B:425:HOH:O	1.80	0.82
1:B:143[B]:LYS:CE	1:B:173:VAL:HG11	2.09	0.81
1:A:80:GLU:O	1:A:84[B]:ILE:HD13	1.83	0.79
1:A:96[B]:ARG:CD	8:A:551:HOH:O	2.30	0.79
1:A:27:VAL:CG2	8:A:656:HOH:O	2.29	0.78
1:A:134:LEU:HD13	8:A:643:HOH:O	1.84	0.78
1:A:80:GLU:O	1:A:84[B]:ILE:CD1	2.33	0.77
1:B:62:TYR:CE2	1:B:66:LEU:CD1	2.68	0.76
1:A:79:VAL:CG2	1:A:82[A]:ARG:HH21	1.98	0.76
1:A:134:LEU:HD22	1:A:143:LYS:NZ	2.01	0.76
1:B:173:VAL:HG13	1:B:178:ALA:HB3	1.68	0.75
1:B:114[A]:LEU:HD21	1:B:217[A]:MET:SD	2.26	0.75
1:A:134:LEU:HD22	1:A:143:LYS:HZ3	1.49	0.75
1:A:135:TRP:CB	5:A:305:EDO:H22	2.17	0.75
1:B:84[B]:ILE:HD11	1:B:236:VAL:HG11	1.68	0.74
1:A:143:LYS:CE	1:A:173:VAL:HG11	2.17	0.73
1:A:108:HIS:NE2	8:A:654:HOH:O	2.23	0.72
1:A:139:LYS:HE3	8:A:433:HOH:O	1.90	0.71
1:A:180:OAF:HZ3	8:A:692:HOH:O	1.89	0.71
1:A:82[B]:ARG:NH1	8:A:502:HOH:O	2.24	0.70
1:B:122:LYS:CG	8:B:425:HOH:O	2.38	0.70
1:A:139:LYS:CE	8:A:433:HOH:O	2.39	0.69
1:B:80:GLU:O	1:B:84[B]:ILE:CD1	2.40	0.69
1:A:79:VAL:HG23	1:A:82[A]:ARG:HH21	1.57	0.68
1:B:160:ASN:HB2	1:B:161:PRO:HD2	1.75	0.68
1:B:226:ILE:HG21	8:B:627:HOH:O	1.93	0.68
1:A:135:TRP:HB2	5:A:305:EDO:H22	1.76	0.67
1:A:135:TRP:H	5:A:305:EDO:C2	2.08	0.67
1:A:16[B]:ASN:ND2	8:A:546:HOH:O	2.28	0.65
1:A:135:TRP:N	5:A:305:EDO:H22	2.11	0.65
1:A:108:HIS:CE1	1:A:114:LEU:HD11	2.32	0.65
1:B:108:HIS:NE2	8:B:646:HOH:O	2.26	0.65
1:B:32:HIS:HB3	8:B:632:HOH:O	1.96	0.64
1:A:135:TRP:H	5:A:305:EDO:C1	2.10	0.64
1:A:204:LEU:N	8:A:692:HOH:O	2.29	0.64
1:A:189:LYS:HB2	8:A:691:HOH:O	1.95	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143[B]:LYS:HD3	1:B:173:VAL:HG21	1.77	0.64
1:B:33:LEU:HG	8:B:642:HOH:O	1.98	0.63
1:A:239:PHE:C	8:A:591:HOH:O	2.40	0.63
1:A:179:GLN:NE2	5:A:306:EDO:C2	2.53	0.62
1:B:108:HIS:CE1	1:B:114[B]:LEU:HG	2.35	0.62
1:A:108:HIS:CE1	8:A:654:HOH:O	2.53	0.61
1:A:74:ARG:HB2	5:A:304:EDO:H22	1.82	0.61
1:B:102[B]:THR:OG1	1:B:127:VAL:HG12	2.00	0.61
1:B:102[B]:THR:HG22	1:B:103:LEU:H	1.67	0.59
1:B:134:LEU:HD22	1:B:143[B]:LYS:NZ	2.17	0.59
1:B:122:LYS:O	1:B:122:LYS:CG	2.38	0.57
1:B:102[B]:THR:HG22	1:B:103:LEU:N	2.18	0.57
1:B:32:HIS:O	1:B:36:GLU:HG3	2.04	0.57
1:A:63:GLN:CB	8:A:656:HOH:O	2.36	0.57
1:A:75:ASP:OD1	5:A:304:EDO:O2	2.21	0.56
1:A:11[B]:GLU:HG2	8:A:517:HOH:O	2.07	0.55
1:B:48:LYS:HG3	1:B:51[B]:ARG:NH2	2.21	0.55
1:B:239:PHE:C	8:B:556:HOH:O	2.49	0.55
1:A:11[B]:GLU:HG3	1:A:45:VAL:CG2	2.37	0.55
1:A:80:GLU:O	1:A:84[B]:ILE:HD12	2.06	0.55
1:A:135:TRP:N	5:A:305:EDO:C2	2.69	0.54
1:B:114[A]:LEU:HD11	8:B:417:HOH:O	2.06	0.54
1:A:135:TRP:H	5:A:305:EDO:H22	1.65	0.54
1:A:160:ASN:HB2	1:A:161:PRO:CD	2.38	0.54
1:A:138[B]:ASN:ND2	8:A:422:HOH:O	2.22	0.53
1:B:202[B]:LYS:HE3	8:B:414:HOH:O	2.08	0.53
1:B:108:HIS:CG	1:B:114[B]:LEU:HG	2.44	0.53
1:B:134:LEU:HD22	1:B:143[B]:LYS:HZ2	1.74	0.53
1:A:154[B]:VAL:HG22	1:A:178:ALA:HB2	1.90	0.53
1:B:110:ASN:O	1:B:110:ASN:ND2	2.43	0.52
1:A:139:LYS:HD2	8:A:630:HOH:O	2.10	0.51
1:A:154[B]:VAL:HG12	1:A:155:LEU:O	2.10	0.51
1:A:186[B]:LYS:HG3	1:A:190:ASP:O	2.11	0.51
7:B:305:PEG:H12	8:B:476:HOH:O	2.10	0.51
1:A:108:HIS:CE1	1:A:114:LEU:CD1	2.94	0.51
1:A:154[B]:VAL:HG12	1:A:155:LEU:N	2.25	0.50
1:A:11[A]:GLU:HG3	1:A:41:VAL:HG13	1.93	0.50
1:A:79:VAL:CG2	1:A:82[A]:ARG:NH2	2.72	0.50
1:B:139[A]:LYS:HE2	8:B:608:HOH:O	2.11	0.50
1:A:104:SER:O	1:A:113:ARG:HD3	2.13	0.49
1:B:84[B]:ILE:CD1	1:B:236:VAL:HG11	2.39	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:SER:O	1:B:113:ARG:HD3	2.13	0.49
1:B:55[A]:MET:HG2	8:B:447:HOH:O	2.12	0.49
1:B:108:HIS:CG	1:B:114[B]:LEU:CD2	2.96	0.49
1:A:151:GLN:HE21	1:A:171:ASN:ND2	2.12	0.48
1:A:135:TRP:H	5:A:305:EDO:H12	1.77	0.48
1:B:55[A]:MET:HB3	1:B:55[A]:MET:HE3	1.36	0.48
1:A:143:LYS:HD3	1:A:173:VAL:HG21	1.95	0.48
1:B:231:HIS:HD2	8:B:561:HOH:O	1.96	0.47
1:A:80:GLU:CD	8:A:614:HOH:O	2.57	0.47
1:A:108:HIS:CD2	1:A:114:LEU:HG	2.49	0.47
1:A:154[B]:VAL:HG22	1:A:178:ALA:CB	2.45	0.47
1:B:79[A]:VAL:HG13	1:B:82:ARG:NH2	2.31	0.46
1:A:175:SER:HB3	8:A:634:HOH:O	2.15	0.46
1:B:139[B]:LYS:HD3	8:B:624:HOH:O	2.14	0.46
1:A:11[B]:GLU:HG3	1:A:45:VAL:HG23	1.97	0.46
1:A:80:GLU:OE2	8:A:614:HOH:O	2.20	0.45
1:B:108:HIS:ND1	1:B:114[B]:LEU:HG	2.32	0.45
1:A:27:VAL:HG21	8:A:656:HOH:O	2.05	0.45
1:B:95:LYS:NZ	8:B:623:HOH:O	2.50	0.45
1:A:124:SER:O	1:A:127:VAL:HG22	2.16	0.45
1:B:90:ILE:HD12	1:B:90:ILE:C	2.42	0.45
1:A:96[A]:ARG:CZ	1:A:231[A]:HIS:CE1	3.00	0.44
1:A:122:LYS:HD3	8:A:619:HOH:O	2.17	0.44
1:B:166[B]:MET:CE	1:B:219:TRP:CH2	3.00	0.44
1:A:189:LYS:CB	8:A:691:HOH:O	2.61	0.44
1:B:114[A]:LEU:CD2	1:B:217[A]:MET:SD	3.03	0.44
1:B:186:LYS:HD2	1:B:186:LYS:HA	1.88	0.44
1:A:108:HIS:CE1	1:A:114:LEU:HG	2.53	0.44
1:B:16:ASN:CB	8:B:617:HOH:O	2.41	0.43
1:A:173:VAL:HG13	1:A:178:ALA:HB3	1.98	0.43
1:A:143:LYS:CD	1:A:173:VAL:HG21	2.48	0.43
1:A:135:TRP:CZ2	1:A:138[B]:ASN:HA	2.54	0.43
1:B:202[B]:LYS:HE3	8:B:453:HOH:O	2.18	0.43
1:B:180:OAF:HZ3	8:B:649:HOH:O	2.18	0.43
1:B:217[B]:MET:HB2	1:B:217[B]:MET:HE3	1.53	0.43
1:A:28:GLU:HG2	1:A:32:HIS:CE1	2.54	0.42
1:A:237:LYS:HD3	8:A:623:HOH:O	2.19	0.42
1:A:11[B]:GLU:HG3	1:A:45:VAL:HG22	2.00	0.42
1:B:21:ALA:HB2	8:B:609:HOH:O	2.04	0.41
1:A:111:ASP:HB3	1:A:168:PHE:CE1	2.55	0.41
1:A:189:LYS:C	8:A:691:HOH:O	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96[B]:ARG:HD2	8:A:551:HOH:O	2.10	0.41
1:B:108:HIS:CG	1:B:114[B]:LEU:HD21	2.56	0.41
1:A:108:HIS:NE2	1:A:114:LEU:HG	2.37	0.40
1:B:160:ASN:HB2	1:B:161:PRO:CD	2.48	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	248/239 (104%)	242 (98%)	6 (2%)	0	100	100
1	B	247/239 (103%)	242 (98%)	5 (2%)	0	100	100
All	All	495/478 (104%)	484 (98%)	11 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	220/207 (106%)	219 (100%)	1 (0%)	81	61
1	B	219/207 (106%)	214 (98%)	5 (2%)	44	11
All	All	439/414 (106%)	433 (99%)	6 (1%)	63	24

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	LEU
1	B	55[A]	MET
1	B	55[B]	MET
1	B	120	LYS
1	B	122	LYS
1	B	134	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	52	ASN
1	A	67	GLN
1	A	150	ASN
1	A	171	ASN
1	A	179	GLN
1	B	53	ASN
1	B	231	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	0AF	B	180	1	15,16,17	3.78	3 (20%)	14,22,24	4.08	8 (57%)
1	0AF	A	180	1	15,16,17	2.82	3 (20%)	14,22,24	3.41	7 (50%)

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
1	0AF	B	180	1	-	0/5/6/8	0/2/2/2
1	0AF	A	180	1	-	0/5/6/8	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	180	0AF	O1-CZ2	-13.33	1.09	1.36
1	A	180	0AF	O1-CZ2	-8.88	1.18	1.36
1	A	180	0AF	CD2-CG	-4.58	1.36	1.44
1	B	180	0AF	CD2-CG	-4.52	1.36	1.44
1	A	180	0AF	CE2-CZ2	-2.75	1.33	1.40
1	B	180	0AF	CE2-CZ2	-2.07	1.35	1.40

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	0AF	O1-CZ2-CE2	-8.63	91.96	118.53
1	B	180	0AF	O1-CZ2-CE2	-7.98	93.96	118.53
1	B	180	0AF	CH2-CZ2-CE2	6.03	129.12	118.69
1	B	180	0AF	CE2-NE1-CD1	-5.87	103.87	108.93
1	B	180	0AF	O1-CZ2-CH2	5.04	132.89	119.36
1	B	180	0AF	CD2-CE2-NE1	4.91	114.52	107.54
1	A	180	0AF	O1-CZ2-CH2	4.83	132.32	119.36
1	B	180	0AF	CZ3-CH2-CZ2	-4.66	114.30	120.05
1	A	180	0AF	CZ3-CH2-CZ2	-4.17	114.89	120.05
1	A	180	0AF	CH2-CZ2-CE2	4.01	125.62	118.69
1	A	180	0AF	CE2-NE1-CD1	-3.83	105.63	108.93
1	B	180	0AF	CD2-CG-CD1	3.80	110.05	106.16
1	A	180	0AF	CD2-CE2-NE1	2.62	111.27	107.54
1	B	180	0AF	CE3-CD2-CE2	2.21	121.62	119.14
1	A	180	0AF	CD2-CG-CD1	2.12	108.33	106.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	180	0AF	1	0
1	A	180	0AF	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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	$\# Z > 2$
2	SCN	A	301	3	1,2,2	1.71	0	0,1,1	-	-
5	EDO	B	304	-	3,3,3	0.41	0	2,2,2	0.52	0
5	EDO	A	306	-	3,3,3	0.48	0	2,2,2	0.30	0
5	EDO	A	305	-	3,3,3	0.52	0	2,2,2	0.42	0
6	PGE	A	307	-	9,9,9	0.35	0	8,8,8	0.90	0
7	PEG	B	305	-	6,6,6	0.37	0	5,5,5	1.20	0
5	EDO	A	304	-	3,3,3	0.41	0	2,2,2	0.76	0
2	SCN	B	301	3	1,2,2	1.05	0	0,1,1	-	-

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
5	EDO	B	304	-	-	1/1/1/1	-
5	EDO	A	306	-	-	1/1/1/1	-
5	EDO	A	305	-	-	0/1/1/1	-
6	PGE	A	307	-	-	3/7/7/7	-
7	PEG	B	305	-	-	2/4/4/4	-
5	EDO	A	304	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	307	PGE	O2-C3-C4-O3
5	B	304	EDO	O1-C1-C2-O2
5	A	306	EDO	O1-C1-C2-O2
5	A	304	EDO	O1-C1-C2-O2
6	A	307	PGE	O3-C5-C6-O4
6	A	307	PGE	C3-C4-O3-C5
7	B	305	PEG	O1-C1-C2-O2
7	B	305	PEG	C1-C2-O2-C3

There are no ring outliers.

4 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	306	EDO	4	0
5	A	305	EDO	8	0
7	B	305	PEG	1	0
5	A	304	EDO	2	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	234/239 (97%)	0.65	17 (7%)	21 30	6, 13, 24, 30	16 (6%)
1	B	234/239 (97%)	0.53	13 (5%)	30 41	6, 13, 23, 29	15 (6%)
All	All	468/478 (97%)	0.59	30 (6%)	25 35	6, 13, 24, 30	31 (6%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	173	VAL	5.1
1	A	239	PHE	4.8
1	A	173	VAL	4.1
1	A	190	ASP	4.0
1	B	108	HIS	3.7
1	A	212	THR	3.7
1	B	109	GLY	3.3
1	A	172	SER	3.3
1	A	32	HIS	3.2
1	A	174	ASP	3.1
1	A	108	HIS	2.9
1	B	172	SER	2.8
1	A	110	ASN	2.8
1	B	190	ASP	2.7
1	B	212	THR	2.6
1	A	5	VAL	2.6
1	A	123	THR	2.5
1	A	188	ASP	2.5
1	B	8	ASP	2.4
1	A	226	ILE	2.3
1	A	120	LYS	2.3
1	B	110	ASN	2.3
1	A	187	TYR	2.3
1	B	5	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	186[A]	LYS	2.2
1	B	189	LYS	2.2
1	B	162	ASN	2.1
1	B	32	HIS	2.1
1	B	239	PHE	2.0
1	A	119	GLY	2.0

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

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. 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	B-factors(Å ²)	Q<0.9
1	0AF	A	180	15/16	0.96	0.06	6,8,10,10	0
1	0AF	B	180	15/16	0.96	0.06	7,8,11,11	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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. 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	B-factors(Å ²)	Q<0.9
5	EDO	B	304	4/4	0.76	0.14	32,32,32,34	0
7	PEG	B	305	7/7	0.81	0.14	23,25,29,29	0
5	EDO	A	306	4/4	0.85	0.12	22,25,29,31	0
6	PGE	A	307	10/10	0.88	0.10	24,28,30,31	0
5	EDO	A	304	4/4	0.88	0.12	25,27,29,31	0
5	EDO	A	305	4/4	0.93	0.10	17,20,23,24	0
4	K	A	303	1/1	0.98	0.04	20,20,20,20	0
4	K	B	303	1/1	0.98	0.08	31,31,31,31	0
2	SCN	A	301	3/3	0.98	0.07	10,10,11,14	0
2	SCN	B	301	3/3	0.98	0.07	10,10,11,14	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CD	A	302	1/1	1.00	0.01	10,10,10,10	0
3	CD	B	302	1/1	1.00	0.01	10,10,10,10	0

6.5 Other polymers [i](#)

There are no such residues in this entry.