



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:58 PM UTC

PDB ID : 4A34 / pdb_00004a34
Title : Crystal structure of the fucose mutarotase in complex with L-fucose from *Streptococcus pneumoniae*
Authors : Higgins, M.A.; Boraston, A.B.
Deposited on : 2011-09-29
Resolution : 2.50 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

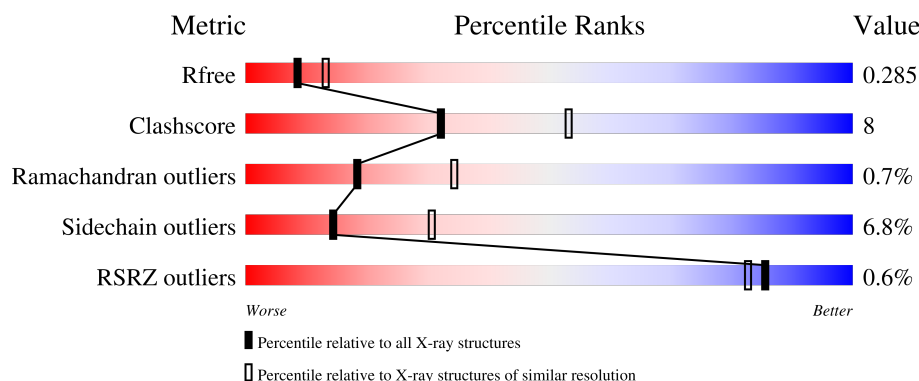
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	147	% 85% 9% • 5%
1	B	147	% 74% 17% • 5%
1	C	147	% 78% 14% • 5%
1	D	147	% 80% 14% • 5%
1	E	147	% 73% 20% • 5%

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Mol	Chain	Length	Quality of chain
1	F	147	
1	G	147	
1	H	147	
1	I	147	
1	J	147	
1	K	147	
1	L	147	
1	M	147	
1	N	147	
1	O	147	
1	P	147	
1	Q	147	
1	R	147	
1	S	147	
1	T	147	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 22775 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 RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	140	Total	C	N	O	S	0	0	0
			1087	694	178	208	7			
1	B	140	Total	C	N	O	S	0	0	0
			1080	690	178	205	7			
1	C	140	Total	C	N	O	S	0	0	0
			1073	685	177	204	7			
1	D	140	Total	C	N	O	S	0	1	0
			1095	699	178	211	7			
1	E	139	Total	C	N	O	S	0	0	0
			1075	685	175	208	7			
1	F	140	Total	C	N	O	S	0	0	0
			1073	686	176	204	7			
1	G	140	Total	C	N	O	S	0	0	0
			1098	702	180	209	7			
1	H	140	Total	C	N	O	S	0	0	0
			1089	696	178	208	7			
1	I	142	Total	C	N	O	S	0	0	0
			1100	702	180	211	7			
1	J	140	Total	C	N	O	S	0	0	0
			1076	687	177	205	7			
1	K	140	Total	C	N	O	S	0	0	0
			1086	693	177	209	7			
1	L	140	Total	C	N	O	S	0	0	0
			1098	702	180	209	7			
1	M	140	Total	C	N	O	S	0	0	0
			1094	699	179	209	7			
1	N	140	Total	C	N	O	S	0	0	0
			1075	689	174	205	7			
1	O	140	Total	C	N	O	S	0	0	0
			1079	689	177	206	7			
1	P	140	Total	C	N	O	S	0	0	0
			1080	689	177	207	7			

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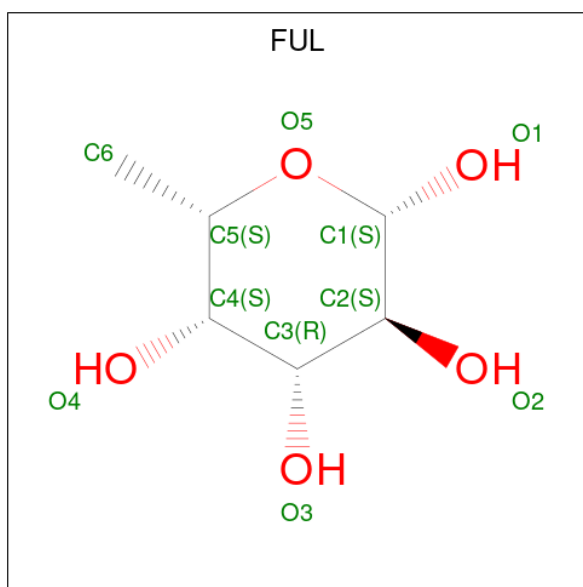
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	140	Total	C	N	O	S	0	0	0
			1084	692	177	208	7			
1	R	140	Total	C	N	O	S	0	0	0
			1080	691	175	207	7			
1	S	140	Total	C	N	O	S	0	0	0
			1094	702	180	205	7			
1	T	140	Total	C	N	O	S	0	0	0
			1079	690	177	205	7			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	147	GLN	-	expression tag	UNP Q04I09
B	147	GLN	-	expression tag	UNP Q04I09
C	147	GLN	-	expression tag	UNP Q04I09
D	147	GLN	-	expression tag	UNP Q04I09
E	147	GLN	-	expression tag	UNP Q04I09
F	147	GLN	-	expression tag	UNP Q04I09
G	147	GLN	-	expression tag	UNP Q04I09
H	147	GLN	-	expression tag	UNP Q04I09
I	147	GLN	-	expression tag	UNP Q04I09
J	147	GLN	-	expression tag	UNP Q04I09
K	147	GLN	-	expression tag	UNP Q04I09
L	147	GLN	-	expression tag	UNP Q04I09
M	147	GLN	-	expression tag	UNP Q04I09
N	147	GLN	-	expression tag	UNP Q04I09
O	147	GLN	-	expression tag	UNP Q04I09
P	147	GLN	-	expression tag	UNP Q04I09
Q	147	GLN	-	expression tag	UNP Q04I09
R	147	GLN	-	expression tag	UNP Q04I09
S	147	GLN	-	expression tag	UNP Q04I09
T	147	GLN	-	expression tag	UNP Q04I09

- Molecule 2 is beta-L-fucopyranose (CCD ID: FUL) (formula: C₆H₁₂O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			11	6	5		
2	B	1	Total	C	O	0	0
			11	6	5		
2	C	1	Total	C	O	0	0
			11	6	5		
2	D	1	Total	C	O	0	0
			11	6	5		
2	E	1	Total	C	O	0	0
			11	6	5		
2	F	1	Total	C	O	0	0
			11	6	5		
2	G	1	Total	C	O	0	0
			11	6	5		
2	H	1	Total	C	O	0	0
			11	6	5		
2	I	1	Total	C	O	0	0
			11	6	5		
2	J	1	Total	C	O	0	0
			11	6	5		
2	K	1	Total	C	O	0	0
			11	6	5		
2	L	1	Total	C	O	0	0
			11	6	5		
2	M	1	Total	C	O	0	0
			11	6	5		
2	N	1	Total	C	O	0	0
			11	6	5		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	O	1	Total C O 11 6 5	0	0
2	P	1	Total C O 11 6 5	0	0
2	Q	1	Total C O 11 6 5	0	0
2	R	1	Total C O 11 6 5	0	0
2	S	1	Total C O 11 6 5	0	0
2	T	1	Total C O 11 6 5	0	0

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total K 1 1	0	0
3	C	1	Total K 1 1	0	0
3	D	1	Total K 1 1	0	0
3	E	1	Total K 1 1	0	0
3	J	1	Total K 1 1	0	0
3	K	1	Total K 1 1	0	0
3	L	1	Total K 1 1	0	0
3	N	1	Total K 1 1	0	0
3	Q	1	Total K 1 1	0	0
3	S	1	Total K 1 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	58	Total O 58 58	0	0

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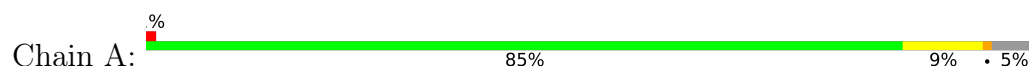
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	49	Total 49	O 49	0	0
4	C	46	Total 46	O 46	0	0
4	D	38	Total 38	O 38	0	0
4	E	58	Total 58	O 58	0	0
4	F	44	Total 44	O 44	0	0
4	G	28	Total 28	O 28	0	0
4	H	41	Total 41	O 41	0	0
4	I	30	Total 30	O 30	0	0
4	J	22	Total 22	O 22	0	0
4	K	56	Total 56	O 56	0	0
4	L	31	Total 31	O 31	0	0
4	M	46	Total 46	O 46	0	0
4	N	50	Total 50	O 50	0	0
4	O	44	Total 44	O 44	0	0
4	P	55	Total 55	O 55	0	0
4	Q	27	Total 27	O 27	0	0
4	R	34	Total 34	O 34	0	0
4	S	48	Total 48	O 48	0	0
4	T	45	Total 45	O 45	0	0

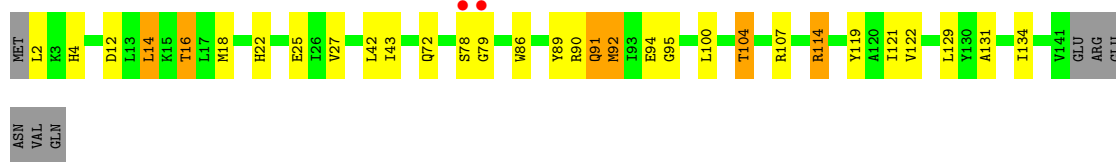
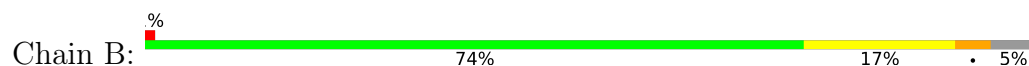
3 Residue-property plots [i](#)

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.

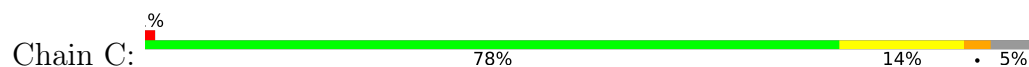
• Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN



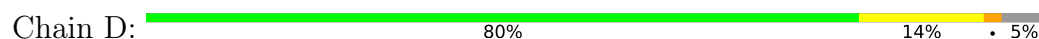
• Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN



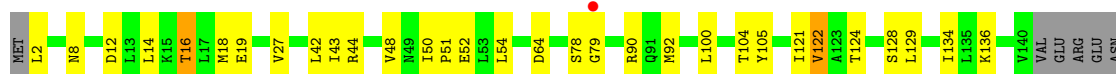
• Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN



• Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN



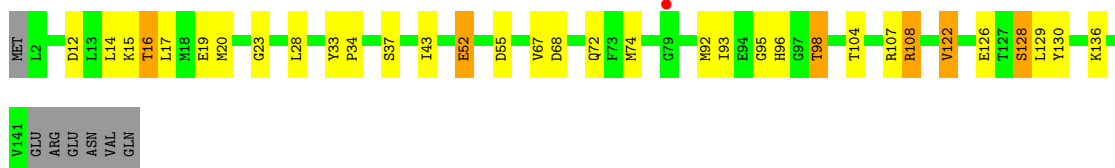
• Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN



VAL
GLN

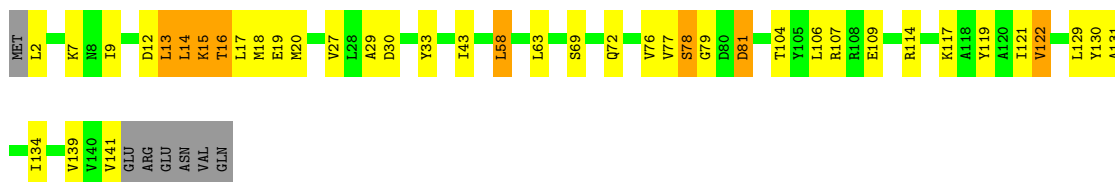
- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain F:  %



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain G:  %



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain H:  %



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain I:  %



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain J:  %



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain K:  %



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain L: 78% 14% 5%



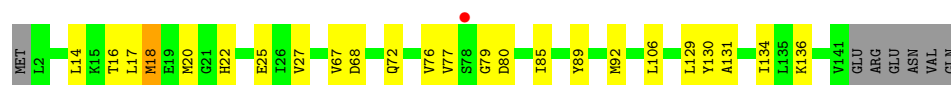
- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain M: 77% 17% 5%



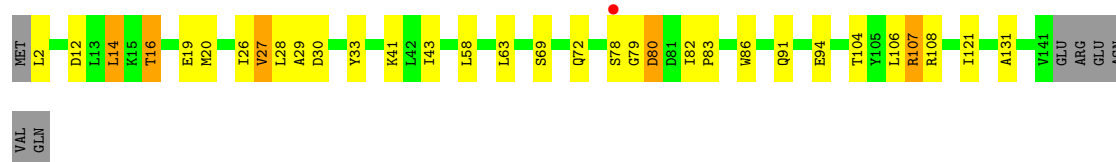
- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain N: 79% 16% 5%



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain O: 73% 18% 5%



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain P: 76% 17% 5%



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

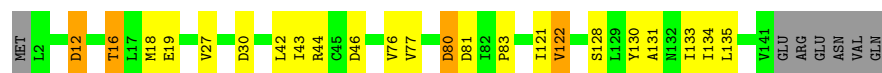
Chain Q: 66% 25% 5%





- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain R: 80% 13% • 5%



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain S: 77% 14% • 5%



- Molecule 1: RBSD/FUCU TRANSPORT PROTEIN FAMILY PROTEIN

Chain T: 78% 14% • • 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.32Å 144.16Å 165.46Å 90.00° 97.51° 90.00°	Depositor
Resolution (Å)	19.97 – 2.50 19.97 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.97-2.50) 99.4 (19.97-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.50Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.225 , 0.292 0.220 , 0.285	Depositor DCC
R_{free} test set	5139 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.042	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	22775	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.10% 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: K, FUL

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	0.53	0/1107	0.83	1/1502 (0.1%)
1	B	0.53	0/1100	0.85	0/1492
1	C	0.56	0/1093	0.82	0/1485
1	D	0.53	0/1115	0.81	0/1513
1	E	0.56	0/1095	0.84	0/1488
1	F	0.54	0/1093	0.86	0/1485
1	G	0.57	0/1118	0.85	0/1514
1	H	0.52	0/1109	0.82	0/1505
1	I	0.55	0/1120	0.85	0/1520
1	J	0.56	0/1096	0.84	0/1489
1	K	0.53	0/1106	0.86	0/1502
1	L	0.51	0/1118	0.83	1/1514 (0.1%)
1	M	0.55	0/1114	0.86	0/1510
1	N	0.56	0/1095	0.87	0/1488
1	O	0.54	0/1099	0.83	0/1493
1	P	0.56	0/1100	0.85	2/1494 (0.1%)
1	Q	0.53	0/1104	0.83	0/1498
1	R	0.53	0/1100	0.82	0/1494
1	S	0.56	0/1114	0.85	1/1506 (0.1%)
1	T	0.57	0/1099	0.82	0/1493
All	All	0.55	0/22095	0.84	5/29985 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	81	ASP	N-CA-C	6.60	118.47	111.28
1	P	95	GLY	N-CA-C	-5.56	107.63	113.58
1	A	39	ALA	N-CA-C	5.32	117.41	109.59
1	P	77	VAL	N-CA-C	5.26	115.41	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	S	40	ASN	N-CA-C	-5.15	106.16	112.90

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	1087	0	1084	10	0
1	B	1080	0	1076	20	0
1	C	1073	0	1058	18	0
1	D	1095	0	1093	17	0
1	E	1075	0	1063	21	0
1	F	1073	0	1064	25	0
1	G	1098	0	1116	31	0
1	H	1089	0	1091	14	0
1	I	1100	0	1098	18	0
1	J	1076	0	1068	18	0
1	K	1086	0	1083	18	0
1	L	1098	0	1116	15	0
1	M	1094	0	1105	14	0
1	N	1075	0	1070	18	0
1	O	1079	0	1069	26	0
1	P	1080	0	1072	24	0
1	Q	1084	0	1081	25	0
1	R	1080	0	1079	12	0
1	S	1094	0	1116	17	0
1	T	1079	0	1077	20	0
2	A	11	0	12	0	0
2	B	11	0	12	0	0
2	C	11	0	12	1	0
2	D	11	0	12	0	0
2	E	11	0	12	1	0
2	F	11	0	12	3	0
2	G	11	0	12	0	0
2	H	11	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	I	11	0	12	2	0
2	J	11	0	12	0	0
2	K	11	0	12	0	0
2	L	11	0	12	0	0
2	M	11	0	12	0	0
2	N	11	0	12	0	0
2	O	11	0	12	0	0
2	P	11	0	12	0	0
2	Q	11	0	12	0	0
2	R	11	0	12	0	0
2	S	11	0	12	0	0
2	T	11	0	12	1	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	J	1	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
3	N	1	0	0	0	0
3	Q	1	0	0	0	0
3	S	1	0	0	0	0
4	A	58	0	0	0	0
4	B	49	0	0	3	0
4	C	46	0	0	2	0
4	D	38	0	0	1	0
4	E	58	0	0	1	0
4	F	44	0	0	1	0
4	G	28	0	0	1	0
4	H	41	0	0	0	0
4	I	30	0	0	0	0
4	J	22	0	0	0	0
4	K	56	0	0	0	0
4	L	31	0	0	1	0
4	M	46	0	0	0	0
4	N	50	0	0	0	0
4	O	44	0	0	5	0
4	P	55	0	0	0	0
4	Q	27	0	0	1	0
4	R	34	0	0	0	0
4	S	48	0	0	0	0
4	T	45	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	22775	0	21919	351	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 8.

All (351) 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:S:80:ASP:CB	1:S:81:ASP:HA	1.71	1.19
1:S:80:ASP:HB3	1:S:81:ASP:HA	1.20	1.18
1:O:79:GLY:HA2	1:O:80:ASP:HB2	1.22	1.11
1:J:96:HIS:H	1:J:97:GLY:CA	1.63	1.10
1:J:96:HIS:N	1:J:97:GLY:HA2	1.65	1.09
1:K:17:LEU:HA	1:K:20:MET:HE2	1.41	0.99
1:J:96:HIS:H	1:J:97:GLY:HA2	0.82	0.95
1:F:17:LEU:HA	1:F:20:MET:HE3	1.48	0.94
1:S:80:ASP:HB3	1:S:81:ASP:CA	2.01	0.89
1:B:16:THR:HG21	1:B:43:ILE:HD11	1.55	0.88
1:I:16:THR:HG21	1:I:43:ILE:HD11	1.56	0.87
1:G:17:LEU:HA	1:G:20:MET:HE3	1.57	0.86
1:P:12:ASP:O	1:P:16:THR:HG22	1.76	0.85
1:O:79:GLY:CA	1:O:80:ASP:HB2	2.07	0.84
1:O:69:SER:OG	1:O:104:THR:HG22	1.76	0.84
1:O:79:GLY:HA2	1:O:80:ASP:CB	2.04	0.83
1:C:71:ILE:HD12	1:C:93:ILE:HD13	1.57	0.83
1:N:17:LEU:HA	1:N:20:MET:HE3	1.58	0.82
1:P:79:GLY:HA3	1:P:80:ASP:HB2	1.62	0.81
1:K:79:GLY:HA2	1:K:80:ASP:O	1.80	0.81
1:D:16:THR:HG21	1:D:43:ILE:HD11	1.63	0.81
1:S:80:ASP:HB2	1:S:81:ASP:HA	1.64	0.80
1:J:16:THR:HG21	1:J:43:ILE:HD11	1.63	0.80
1:B:2:LEU:HD22	1:G:18:MET:HG2	1.62	0.80
1:P:78:SER:H	1:P:79:GLY:HA2	1.47	0.79
1:O:14:LEU:HD13	1:T:14:LEU:HD13	1.64	0.78
1:J:12:ASP:O	1:J:16:THR:HG22	1.85	0.76
1:G:12:ASP:O	1:G:16:THR:HG23	1.85	0.76
1:N:79:GLY:HA2	1:N:80:ASP:HB2	1.68	0.75
1:E:12:ASP:O	1:E:16:THR:HG23	1.87	0.74
1:Q:94:GLU:HB2	1:Q:100:LEU:HD13	1.69	0.73
1:O:16:THR:HG21	1:O:43:ILE:HD11	1.69	0.72
1:P:79:GLY:CA	1:P:80:ASP:HB2	2.19	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:17:LEU:HA	1:M:20:MET:HE3	1.72	0.71
1:F:12:ASP:O	1:F:16:THR:HG22	1.90	0.71
1:B:12:ASP:O	1:B:16:THR:HG23	1.91	0.70
1:O:12:ASP:O	1:O:16:THR:CG2	2.40	0.70
1:H:141:VAL:HG12	1:H:141:VAL:O	1.90	0.70
1:M:12:ASP:O	1:M:16:THR:HG23	1.90	0.70
1:T:80:ASP:C	1:T:82:ILE:H	1.98	0.70
1:L:14:LEU:HD12	1:Q:18:MET:SD	2.32	0.70
1:O:58:LEU:HD22	1:O:63:LEU:HD11	1.74	0.70
1:Q:27:VAL:HG13	1:Q:121:ILE:HG12	1.74	0.68
1:B:27:VAL:HG22	1:B:134:ILE:HG12	1.75	0.68
1:F:72:GLN:HG3	1:F:104:THR:HG22	1.75	0.68
1:A:72:GLN:HG3	1:A:104:THR:CG2	2.24	0.67
1:S:80:ASP:CB	1:S:81:ASP:CA	2.56	0.67
1:F:74:MET:HE3	1:F:130:TYR:CD1	2.28	0.67
1:H:78:SER:CB	1:H:79:GLY:CA	2.72	0.67
1:P:16:THR:HB	1:P:41:LYS:NZ	2.10	0.66
1:O:12:ASP:O	1:O:16:THR:HG22	1.94	0.66
1:O:20:MET:HE1	1:O:26:ILE:HD12	1.77	0.66
1:R:12:ASP:O	1:R:16:THR:HG22	1.95	0.66
1:H:76:VAL:HA	1:H:130:TYR:CE1	2.31	0.66
1:E:18:MET:SD	1:J:14:LEU:HD23	2.36	0.65
1:F:52:GLU:HG3	4:F:2024:HOH:O	1.96	0.65
1:F:17:LEU:HA	1:F:20:MET:CE	2.24	0.65
1:J:96:HIS:N	1:J:97:GLY:CA	2.39	0.65
1:K:77:VAL:O	1:K:80:ASP:HB2	1.96	0.65
1:A:69:SER:OG	1:A:104:THR:HG22	1.97	0.64
1:B:72:GLN:HG3	1:B:104:THR:HG22	1.78	0.64
1:C:14:LEU:HD12	1:H:18:MET:SD	2.36	0.64
1:G:72:GLN:HG3	1:G:104:THR:CG2	2.28	0.64
1:I:12:ASP:O	1:I:16:THR:HG22	1.97	0.64
1:A:9:ILE:HD11	1:A:14:LEU:HB2	1.79	0.64
1:P:25:GLU:OE2	1:P:136:LYS:HE2	1.97	0.64
1:K:25:GLU:OE1	1:K:114:ARG:HD3	1.98	0.63
1:C:14:LEU:HD13	1:H:14:LEU:HD13	1.81	0.63
1:L:13:LEU:O	1:L:17:LEU:HG	1.99	0.63
1:H:78:SER:CB	1:H:79:GLY:HA3	2.29	0.63
1:S:86:TRP:O	1:S:90:ARG:HG3	1.99	0.62
1:P:78:SER:N	1:P:79:GLY:HA2	2.10	0.62
1:A:72:GLN:HG3	1:A:104:THR:HG23	1.81	0.62
1:G:117:LYS:HG3	4:G:2026:HOH:O	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:16:THR:HG21	1:E:43:ILE:HD11	1.80	0.62
1:F:74:MET:HE1	2:F:1142:FUL:H3	1.81	0.61
1:K:12:ASP:O	1:K:16:THR:CG2	2.49	0.61
1:G:17:LEU:HA	1:G:20:MET:CE	2.28	0.61
1:N:22:HIS:HE1	1:O:108:ARG:HH22	1.49	0.61
1:P:43:ILE:O	1:P:122:VAL:HA	2.01	0.61
1:N:16:THR:HG22	1:N:20:MET:HE2	1.82	0.61
1:F:12:ASP:O	1:F:16:THR:CG2	2.49	0.60
1:M:16:THR:HG21	1:M:43:ILE:HD11	1.82	0.60
1:S:25:GLU:OE1	1:S:114:ARG:HD3	2.02	0.60
1:C:80:ASP:O	4:C:2036:HOH:O	2.16	0.59
1:H:49:ASN:HB2	1:H:52:GLU:OE2	2.02	0.59
1:J:16:THR:HG21	1:J:43:ILE:CD1	2.32	0.59
1:B:4:HIS:ND1	1:S:94:GLU:HG3	2.18	0.59
1:R:12:ASP:O	1:R:16:THR:CG2	2.50	0.59
1:J:93:ILE:HG23	1:J:98:THR:HG22	1.84	0.58
1:J:27:VAL:HG22	1:J:134:ILE:HG12	1.86	0.58
1:T:78:SER:H	1:T:79:GLY:HA3	1.67	0.58
1:P:79:GLY:HA3	1:P:80:ASP:O	2.04	0.57
1:Q:140:VAL:O	1:Q:141:VAL:HB	2.04	0.57
1:G:72:GLN:HG3	1:G:104:THR:HG23	1.86	0.57
1:Q:76:VAL:HG13	1:Q:80:ASP:HB3	1.86	0.57
1:E:27:VAL:HG22	1:E:134:ILE:HG12	1.87	0.57
1:N:22:HIS:CE1	1:O:108:ARG:HH12	2.22	0.56
1:P:16:THR:HG21	1:P:43:ILE:HD11	1.88	0.56
1:B:91:GLN:H	1:B:94:GLU:H	1.51	0.56
1:F:95:GLY:O	1:F:96:HIS:HB2	2.05	0.56
1:L:27:VAL:HG22	1:L:134:ILE:HG12	1.87	0.56
1:O:14:LEU:CD1	1:T:14:LEU:HD13	2.35	0.56
4:O:2031:HOH:O	1:T:4:HIS:HE1	1.88	0.56
1:Q:82:ILE:HD12	1:Q:82:ILE:H	1.70	0.56
1:D:25:GLU:OE1	1:D:114:ARG:HD3	2.06	0.56
1:D:15:LYS:HD2	1:I:15:LYS:HD2	1.88	0.55
1:N:77:VAL:O	1:N:80:ASP:HB2	2.06	0.55
1:E:14:LEU:O	1:E:18:MET:HG3	2.06	0.55
1:O:27:VAL:HG13	1:O:121:ILE:HG12	1.87	0.55
1:F:67:VAL:HG22	1:F:68:ASP:H	1.72	0.55
1:M:25:GLU:OE2	1:M:136:LYS:HE2	2.08	0.55
1:Q:29:ALA:HB1	1:Q:33:TYR:HB3	1.89	0.55
1:F:17:LEU:HD23	1:F:20:MET:HE1	1.89	0.54
1:O:91:GLN:O	1:O:94:GLU:O	2.24	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:67:VAL:HG22	1:N:68:ASP:H	1.73	0.54
1:N:17:LEU:HD23	1:N:20:MET:HE1	1.90	0.54
1:F:126:GLU:OE2	1:F:128:SER:HB3	2.07	0.54
1:C:18:MET:SD	1:H:14:LEU:HD12	2.48	0.54
1:L:52:GLU:H	1:L:52:GLU:CD	2.16	0.54
1:T:27:VAL:HG22	1:T:134:ILE:HG12	1.90	0.54
1:D:12:ASP:O	1:D:16:THR:HG22	2.08	0.53
1:N:72:GLN:HG2	1:N:106:LEU:CD2	2.39	0.53
1:P:12:ASP:O	1:P:16:THR:CG2	2.53	0.53
1:C:89:TYR:O	1:C:93:ILE:HG23	2.08	0.53
1:L:43:ILE:O	1:L:122:VAL:HA	2.07	0.53
1:A:109:GLU:CD	1:A:109:GLU:H	2.17	0.53
1:M:14:LEU:HD12	1:R:18:MET:SD	2.49	0.53
1:O:16:THR:HA	1:O:19:GLU:HG2	1.91	0.53
1:H:43:ILE:O	1:H:122:VAL:HA	2.09	0.52
1:D:16:THR:HG21	1:D:43:ILE:CD1	2.36	0.52
1:F:93:ILE:O	1:F:98:THR:HG23	2.10	0.52
1:T:80:ASP:C	1:T:82:ILE:N	2.67	0.52
1:E:51:PRO:HA	1:E:92:MET:HE1	1.92	0.52
1:I:25:GLU:OE1	1:I:114:ARG:HD3	2.09	0.52
1:F:72:GLN:HG3	1:F:104:THR:CG2	2.38	0.52
4:C:2014:HOH:O	1:D:44:ARG:NH2	2.42	0.52
1:I:20:MET:HE1	1:I:26:ILE:HD12	1.92	0.52
1:R:16:THR:HA	1:R:19:GLU:HG2	1.92	0.52
1:I:63:LEU:CD1	1:I:71:ILE:HD11	2.40	0.52
1:B:91:GLN:HA	1:B:92:MET:C	2.36	0.51
1:G:81:ASP:OD1	1:G:81:ASP:N	2.36	0.51
1:E:64:ASP:OD2	1:E:136:LYS:HD3	2.10	0.51
1:F:43:ILE:O	1:F:122:VAL:HA	2.10	0.51
1:N:72:GLN:HG2	1:N:106:LEU:HD21	1.92	0.51
1:O:16:THR:HG21	1:O:43:ILE:CD1	2.40	0.51
1:T:76:VAL:HG12	1:T:79:GLY:HA3	1.92	0.51
1:G:17:LEU:HD23	1:G:20:MET:HE1	1.92	0.51
1:K:44:ARG:HG2	1:K:46:ASP:OD2	2.11	0.51
1:D:29:ALA:HB1	1:D:33:TYR:HB3	1.92	0.51
1:G:43:ILE:O	1:G:122:VAL:HA	2.11	0.51
1:K:18:MET:HG2	1:P:2:LEU:HD12	1.92	0.51
1:L:81:ASP:O	1:L:82:ILE:HB	2.10	0.51
1:P:79:GLY:HA3	1:P:80:ASP:CB	2.31	0.51
1:O:2:LEU:N	4:O:2001:HOH:O	2.44	0.51
1:T:78:SER:N	1:T:79:GLY:HA3	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:78:SER:OG	1:T:79:GLY:HA2	2.10	0.51
1:I:16:THR:HA	1:I:19:GLU:HG2	1.92	0.50
1:K:12:ASP:O	1:K:16:THR:HG23	2.11	0.50
1:K:76:VAL:HA	1:K:130:TYR:CE1	2.46	0.50
1:F:28:LEU:HD23	1:F:122:VAL:HG13	1.94	0.50
1:G:14:LEU:O	1:G:18:MET:HG3	2.12	0.50
1:I:64:ASP:OD2	1:I:136:LYS:HD3	2.12	0.50
1:R:133:ILE:HG22	1:R:135:LEU:HD12	1.92	0.50
1:I:27:VAL:HG22	1:I:134:ILE:HG12	1.94	0.50
1:A:27:VAL:HB	1:A:121:ILE:HG12	1.95	0.49
1:K:16:THR:HG21	1:K:43:ILE:HD11	1.93	0.49
1:P:16:THR:HB	1:P:41:LYS:HZ1	1.77	0.49
1:L:5:ILE:CD1	1:Q:60:LEU:HD22	2.42	0.49
1:M:16:THR:HA	1:M:19:GLU:HG2	1.94	0.49
1:H:25:GLU:OE1	1:H:114:ARG:HD3	2.13	0.49
1:N:14:LEU:HG	1:S:14:LEU:HD13	1.93	0.49
1:N:76:VAL:HA	1:N:130:TYR:CE1	2.48	0.49
1:B:25:GLU:OE2	1:B:114:ARG:NH1	2.46	0.49
1:K:12:ASP:O	1:K:16:THR:HG22	2.12	0.49
1:N:18:MET:SD	1:S:14:LEU:HD12	2.53	0.49
1:D:14:LEU:HD12	1:I:18:MET:HE2	1.94	0.49
1:E:43:ILE:O	1:E:122:VAL:HA	2.13	0.49
1:G:30:ASP:HB3	1:G:131:ALA:HB3	1.94	0.48
1:L:99:ASP:OD2	1:L:101:LYS:HB3	2.13	0.48
1:O:12:ASP:O	1:O:16:THR:HG23	2.13	0.48
1:L:2:LEU:HD12	1:Q:18:MET:HG2	1.96	0.48
1:M:20:MET:HE1	1:M:26:ILE:HD12	1.94	0.48
1:C:71:ILE:CD1	1:C:93:ILE:HD13	2.36	0.48
1:C:9:ILE:HG21	1:C:14:LEU:HG	1.96	0.48
4:O:2031:HOH:O	1:T:3:LYS:HE2	2.14	0.48
1:P:16:THR:HG21	1:P:43:ILE:CD1	2.44	0.48
1:L:78:SER:HA	1:L:79:GLY:HA2	1.63	0.48
1:M:27:VAL:HG21	1:M:115:SER:HB2	1.96	0.48
1:G:69:SER:OG	1:G:104:THR:HG22	2.14	0.48
1:I:107:ARG:O	1:I:108:ARG:C	2.57	0.48
1:G:12:ASP:O	1:G:16:THR:CG2	2.60	0.47
1:C:90:ARG:HG2	1:C:100:LEU:HD22	1.95	0.47
1:F:74:MET:CE	2:F:1142:FUL:H3	2.45	0.47
1:O:14:LEU:HD12	1:T:18:MET:SD	2.54	0.47
1:S:82:ILE:HD12	1:S:82:ILE:H	1.79	0.47
1:B:18:MET:HG3	1:G:2:LEU:HD12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:HIS:ND1	2:C:1142:FUL:O5	2.45	0.47
1:R:30:ASP:HB3	1:R:131:ALA:HB3	1.95	0.47
1:M:94:GLU:HG2	1:M:100:LEU:HD11	1.97	0.47
1:R:27:VAL:HB	1:R:121:ILE:HG12	1.97	0.47
1:C:50:ILE:N	1:C:51:PRO:HD2	2.29	0.47
1:E:2:LEU:N	4:E:2001:HOH:O	2.46	0.47
1:F:23:GLY:HA2	1:F:136:LYS:HE3	1.96	0.47
1:F:55:ASP:OD1	1:F:96:HIS:HE1	1.98	0.47
1:G:15:LYS:HD3	1:G:19:GLU:OE1	2.14	0.47
1:H:141:VAL:O	1:H:141:VAL:CG1	2.59	0.47
1:I:63:LEU:HD13	1:I:71:ILE:HD11	1.97	0.47
1:R:27:VAL:HG22	1:R:134:ILE:HG12	1.97	0.47
1:S:50:ILE:HG13	1:S:126:GLU:HB2	1.95	0.47
1:E:90:ARG:NH2	1:E:105:TYR:OH	2.48	0.47
1:M:50:ILE:N	1:M:51:PRO:HD2	2.30	0.47
1:B:27:VAL:HB	1:B:121:ILE:HG12	1.96	0.46
1:D:12:ASP:O	1:D:16:THR:CG2	2.63	0.46
1:P:89:TYR:OH	1:P:131:ALA:HB1	2.15	0.46
1:F:107:ARG:O	1:F:108:ARG:C	2.57	0.46
1:G:29:ALA:HB1	1:G:33:TYR:HB3	1.96	0.46
1:B:14:LEU:HD12	1:G:18:MET:SD	2.56	0.46
1:L:49:ASN:HB2	1:L:52:GLU:OE2	2.15	0.46
1:J:107:ARG:HB2	1:J:110:ASP:OD2	2.16	0.46
2:I:1144:FUL:H63	1:J:139:VAL:HG21	1.97	0.46
1:K:77:VAL:O	1:K:78:SER:C	2.59	0.46
1:E:27:VAL:HB	1:E:121:ILE:HG12	1.99	0.45
1:F:67:VAL:HG22	1:F:68:ASP:N	2.30	0.45
1:H:25:GLU:OE2	1:H:136:LYS:HE2	2.16	0.45
1:T:25:GLU:OE1	1:T:114:ARG:HD3	2.17	0.45
1:T:67:VAL:HG21	1:T:136:LYS:HD3	1.99	0.45
1:E:54:LEU:HD23	1:E:92:MET:HE3	1.97	0.45
2:I:1144:FUL:O5	1:J:22:HIS:ND1	2.40	0.45
1:N:79:GLY:HA2	1:N:80:ASP:CB	2.42	0.45
1:T:86:TRP:O	1:T:90:ARG:HG2	2.16	0.45
1:O:72:GLN:HG3	1:O:104:THR:CG2	2.47	0.45
4:B:2048:HOH:O	1:G:2:LEU:N	2.48	0.45
1:K:27:VAL:HG22	1:K:134:ILE:HG12	1.98	0.45
1:L:69:SER:N	4:L:2025:HOH:O	2.48	0.45
1:Q:104:THR:HG22	1:Q:105:TYR:N	2.32	0.45
1:O:29:ALA:HB1	1:O:33:TYR:HB3	1.97	0.45
1:D:89:TYR:CZ	1:D:133:ILE:HD11	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:17:LEU:HD23	1:K:20:MET:HE1	1.98	0.45
1:O:83:PRO:HG2	1:O:86:TRP:CE2	2.52	0.45
1:P:16:THR:HB	1:P:41:LYS:HZ2	1.82	0.45
1:J:12:ASP:O	1:J:16:THR:CG2	2.61	0.44
1:Q:78:SER:HA	1:Q:79:GLY:HA2	1.64	0.44
1:C:43:ILE:O	1:C:122:VAL:HA	2.16	0.44
1:D:59:TYR:O	1:I:3:LYS:HE3	2.17	0.44
1:K:15:LYS:HD2	1:P:12:ASP:HA	1.99	0.44
1:N:17:LEU:HD23	1:N:20:MET:CE	2.46	0.44
1:Q:63:LEU:HD12	1:Q:98:THR:HG21	1.99	0.44
1:E:78:SER:HA	1:E:79:GLY:HA2	1.74	0.44
1:G:77:VAL:O	1:G:79:GLY:N	2.50	0.44
1:D:76:VAL:HA	1:D:130:TYR:CE1	2.52	0.44
1:B:78:SER:HA	1:B:79:GLY:HA2	1.73	0.44
1:F:33:TYR:CG	1:F:34:PRO:HD2	2.52	0.44
1:G:16:THR:HA	1:G:19:GLU:HG2	2.00	0.44
1:O:107:ARG:HG2	4:O:2040:HOH:O	2.16	0.44
1:S:3:LYS:HB3	1:S:4:HIS:CD2	2.53	0.44
1:D:16:THR:HA	1:D:19:GLU:HG2	2.00	0.44
1:I:33:TYR:CD1	1:I:34:PRO:HD2	2.53	0.44
1:J:91:GLN:O	1:J:94:GLU:O	2.36	0.44
1:L:14:LEU:HD13	1:Q:14:LEU:CD2	2.47	0.44
1:D:43:ILE:O	1:D:122:VAL:HA	2.17	0.44
1:I:43:ILE:O	1:I:122:VAL:HA	2.17	0.44
1:K:30:ASP:HB3	1:K:131:ALA:HB3	1.98	0.44
1:Q:89:TYR:O	1:Q:92:MET:HB3	2.18	0.44
1:C:77:VAL:C	1:C:79:GLY:HA3	2.43	0.43
1:Q:94:GLU:HB2	1:Q:100:LEU:CD1	2.42	0.43
1:S:79:GLY:O	1:S:80:ASP:O	2.36	0.43
1:F:74:MET:HE1	2:F:1142:FUL:H1	2.00	0.43
1:I:74:MET:SD	1:I:108:ARG:HG3	2.58	0.43
4:B:2011:HOH:O	1:C:44:ARG:NH2	2.49	0.43
1:K:14:LEU:HD23	1:P:18:MET:SD	2.59	0.43
1:P:64:ASP:OD2	1:P:136:LYS:HD3	2.18	0.43
1:S:27:VAL:HG22	1:S:134:ILE:HG12	1.99	0.43
1:J:50:ILE:HG13	1:J:126:GLU:HB2	2.01	0.43
1:B:2:LEU:N	4:B:2001:HOH:O	2.51	0.43
1:F:15:LYS:O	1:F:19:GLU:HG3	2.17	0.43
1:P:13:LEU:HA	1:P:16:THR:HG23	2.00	0.43
1:R:77:VAL:O	1:R:80:ASP:HB2	2.18	0.43
1:T:27:VAL:HB	1:T:121:ILE:HG12	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:82:ILE:HA	1:T:83:PRO:HD3	1.94	0.43
1:O:30:ASP:HB3	1:O:131:ALA:HB3	2.00	0.43
1:T:130:TYR:CZ	2:T:1142:FUL:H62	2.54	0.43
1:A:69:SER:OG	1:A:104:THR:CG2	2.65	0.43
1:A:78:SER:HA	1:A:79:GLY:HA2	1.68	0.43
1:G:27:VAL:HG22	1:G:134:ILE:HG12	2.00	0.43
1:G:107:ARG:HG2	1:G:109:GLU:OE2	2.18	0.43
1:G:76:VAL:HA	1:G:130:TYR:CE1	2.54	0.43
1:M:89:TYR:CE2	1:M:133:ILE:HD11	2.54	0.43
1:P:27:VAL:HG22	1:P:134:ILE:HG12	2.00	0.43
1:Q:117:LYS:O	4:Q:2024:HOH:O	2.21	0.43
1:T:76:VAL:CG1	1:T:77:VAL:N	2.81	0.43
1:N:25:GLU:CG	1:N:136:LYS:HG3	2.49	0.43
1:B:119:TYR:CD1	1:C:37:SER:HB3	2.54	0.43
1:E:48:VAL:O	1:E:124:THR:HB	2.19	0.43
1:O:41:LYS:NZ	4:O:2010:HOH:O	2.51	0.43
1:P:14:LEU:O	1:P:18:MET:HG3	2.19	0.43
1:Q:25:GLU:OE1	1:Q:114:ARG:HD3	2.18	0.43
1:T:78:SER:N	1:T:79:GLY:CA	2.82	0.42
1:D:29:ALA:HB1	1:D:33:TYR:CB	2.48	0.42
1:H:6:PRO:CB	1:H:56:SER:HB2	2.49	0.42
1:I:12:ASP:O	1:I:16:THR:CG2	2.67	0.42
1:K:17:LEU:HD23	1:K:20:MET:CE	2.49	0.42
1:R:43:ILE:O	1:R:122:VAL:HA	2.19	0.42
1:J:61:MET:HA	1:J:62:PRO:HD2	1.89	0.42
1:Q:27:VAL:CG1	1:Q:121:ILE:HG12	2.46	0.42
1:R:83:PRO:HG3	1:R:130:TYR:HB2	2.01	0.42
1:G:17:LEU:HD23	1:G:20:MET:CE	2.49	0.42
1:Q:55:ASP:OD1	1:Q:96:HIS:HE1	2.03	0.42
1:P:79:GLY:CA	1:P:80:ASP:CB	2.89	0.42
1:H:76:VAL:HG12	1:H:130:TYR:CD2	2.55	0.42
1:N:27:VAL:HG22	1:N:134:ILE:HG12	2.00	0.42
1:C:93:ILE:HD12	1:C:98:THR:HB	2.00	0.42
1:Q:48:VAL:HG11	1:Q:53:LEU:HD13	2.01	0.42
1:C:114:ARG:HG2	1:C:134:ILE:HD13	2.01	0.42
1:E:16:THR:HA	1:E:19:GLU:HG2	2.00	0.42
1:S:107:ARG:HG3	1:S:107:ARG:HH11	1.84	0.42
1:B:91:GLN:HA	1:B:95:GLY:H	1.85	0.41
1:G:9:ILE:HG23	1:G:13:LEU:HB3	2.01	0.41
1:C:14:LEU:O	1:C:18:MET:HG3	2.19	0.41
1:B:94:GLU:HG3	1:B:100:LEU:HG	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:109:GLU:CD	1:G:109:GLU:H	2.28	0.41
1:D:9:ILE:HD11	1:D:56:SER:HB3	2.02	0.41
1:M:33:TYR:CG	1:M:34:PRO:HD2	2.56	0.41
1:L:29:ALA:HB1	1:L:33:TYR:HB3	2.01	0.41
1:M:94:GLU:O	1:M:94:GLU:HG3	2.20	0.41
1:Q:43:ILE:O	1:Q:122:VAL:HA	2.20	0.41
1:Q:83:PRO:HG3	1:Q:130:TYR:HB2	2.03	0.41
1:C:85:ILE:HD13	1:C:128:SER:OG	2.21	0.41
1:E:90:ARG:HG3	1:E:100:LEU:HD22	2.02	0.41
1:G:58:LEU:HD12	1:G:58:LEU:HA	1.91	0.41
1:A:9:ILE:HD11	1:A:14:LEU:CB	2.48	0.41
1:E:8:ASN:ND2	1:E:52:GLU:HG2	2.36	0.41
1:E:54:LEU:HD23	1:E:92:MET:CE	2.51	0.41
1:F:37:SER:HB3	1:G:119:TYR:CD1	2.56	0.41
1:G:27:VAL:HB	1:G:121:ILE:HG12	2.03	0.41
1:L:3:LYS:HD3	1:Q:62:PRO:HG3	2.03	0.41
1:R:44:ARG:HD3	1:R:46:ASP:OD2	2.21	0.41
1:E:50:ILE:N	1:E:51:PRO:HD2	2.35	0.40
1:B:86:TRP:O	1:B:90:ARG:HG3	2.21	0.40
1:B:89:TYR:OH	1:B:131:ALA:HB1	2.21	0.40
1:D:22:HIS:ND1	2:E:1141:FUL:O5	2.45	0.40
1:I:16:THR:HG21	1:I:43:ILE:CD1	2.37	0.40
1:G:16:THR:HG21	1:G:43:ILE:HD11	2.04	0.40
1:M:30:ASP:HB3	1:M:131:ALA:HB3	2.03	0.40
1:N:89:TYR:OH	1:N:131:ALA:HB1	2.20	0.40
1:Q:20:MET:HE1	1:Q:26:ILE:HB	2.04	0.40
1:Q:104:THR:CG2	1:Q:105:TYR:N	2.84	0.40
1:E:18:MET:HG2	1:J:2:LEU:HD12	2.03	0.40
1:S:99:ASP:CG	1:S:101:LYS:HD3	2.46	0.40
1:A:29:ALA:HB1	1:A:33:TYR:HB3	2.04	0.40
4:D:2005:HOH:O	1:E:44:ARG:NH2	2.54	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	138/147 (94%)	134 (97%)	4 (3%)	0	100	100
1	B	138/147 (94%)	131 (95%)	6 (4%)	1 (1%)	18	34
1	C	138/147 (94%)	134 (97%)	3 (2%)	1 (1%)	18	34
1	D	139/147 (95%)	131 (94%)	8 (6%)	0	100	100
1	E	137/147 (93%)	131 (96%)	6 (4%)	0	100	100
1	F	138/147 (94%)	131 (95%)	6 (4%)	1 (1%)	18	34
1	G	138/147 (94%)	130 (94%)	7 (5%)	1 (1%)	18	34
1	H	138/147 (94%)	133 (96%)	5 (4%)	0	100	100
1	I	140/147 (95%)	137 (98%)	3 (2%)	0	100	100
1	J	138/147 (94%)	125 (91%)	11 (8%)	2 (1%)	9	17
1	K	138/147 (94%)	132 (96%)	3 (2%)	3 (2%)	5	9
1	L	138/147 (94%)	135 (98%)	2 (1%)	1 (1%)	18	34
1	M	138/147 (94%)	135 (98%)	3 (2%)	0	100	100
1	N	138/147 (94%)	133 (96%)	5 (4%)	0	100	100
1	O	138/147 (94%)	131 (95%)	5 (4%)	2 (1%)	9	17
1	P	138/147 (94%)	131 (95%)	6 (4%)	1 (1%)	18	34
1	Q	138/147 (94%)	132 (96%)	6 (4%)	0	100	100
1	R	138/147 (94%)	133 (96%)	4 (3%)	1 (1%)	18	34
1	S	138/147 (94%)	130 (94%)	6 (4%)	2 (1%)	9	17
1	T	138/147 (94%)	130 (94%)	4 (3%)	4 (3%)	3	5
All	All	2762/2940 (94%)	2639 (96%)	103 (4%)	20 (1%)	18	34

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	J	78	SER
1	K	79	GLY
1	K	80	ASP
1	O	80	ASP
1	S	78	SER
1	S	80	ASP
1	B	91	GLN
1	C	78	SER

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Mol	Chain	Res	Type
1	F	108	ARG
1	K	78	SER
1	T	78	SER
1	T	81	ASP
1	O	78	SER
1	T	79	GLY
1	G	78	SER
1	J	81	ASP
1	L	82	ILE
1	P	80	ASP
1	R	80	ASP
1	T	77	VAL

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	119/130 (92%)	115 (97%)	4 (3%)	32	60
1	B	117/130 (90%)	108 (92%)	9 (8%)	12	25
1	C	115/130 (88%)	108 (94%)	7 (6%)	17	35
1	D	121/130 (93%)	114 (94%)	7 (6%)	18	38
1	E	118/130 (91%)	112 (95%)	6 (5%)	21	43
1	F	116/130 (89%)	108 (93%)	8 (7%)	14	30
1	G	123/130 (95%)	108 (88%)	15 (12%)	5	10
1	H	120/130 (92%)	110 (92%)	10 (8%)	10	22
1	I	121/130 (93%)	113 (93%)	8 (7%)	15	32
1	J	117/130 (90%)	110 (94%)	7 (6%)	17	36
1	K	120/130 (92%)	113 (94%)	7 (6%)	18	38
1	L	123/130 (95%)	114 (93%)	9 (7%)	13	27
1	M	122/130 (94%)	115 (94%)	7 (6%)	18	39
1	N	117/130 (90%)	113 (97%)	4 (3%)	32	60

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	O	117/130 (90%)	110 (94%)	7 (6%)	17	36
1	P	118/130 (91%)	109 (92%)	9 (8%)	12	26
1	Q	119/130 (92%)	106 (89%)	13 (11%)	6	13
1	R	119/130 (92%)	112 (94%)	7 (6%)	18	37
1	S	123/130 (95%)	112 (91%)	11 (9%)	9	20
1	T	118/130 (91%)	110 (93%)	8 (7%)	14	31
All	All	2383/2600 (92%)	2220 (93%)	163 (7%)	14	31

All (163) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ILE
1	A	14	LEU
1	A	44	ARG
1	A	122	VAL
1	B	14	LEU
1	B	16	THR
1	B	42	LEU
1	B	92	MET
1	B	104	THR
1	B	107	ARG
1	B	114	ARG
1	B	122	VAL
1	B	129	LEU
1	C	14	LEU
1	C	85	ILE
1	C	92	MET
1	C	93	ILE
1	C	98	THR
1	C	122	VAL
1	C	129	LEU
1	D	13	LEU
1	D	14	LEU
1	D	16	THR
1	D	42	LEU
1	D	52[A]	GLU
1	D	52[B]	GLU
1	D	122	VAL
1	E	16	THR
1	E	42	LEU

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Mol	Chain	Res	Type
1	E	104	THR
1	E	122	VAL
1	E	128	SER
1	E	129	LEU
1	F	14	LEU
1	F	16	THR
1	F	52	GLU
1	F	92	MET
1	F	98	THR
1	F	122	VAL
1	F	128	SER
1	F	129	LEU
1	G	7	LYS
1	G	13	LEU
1	G	14	LEU
1	G	15	LYS
1	G	16	THR
1	G	58	LEU
1	G	63	LEU
1	G	78	SER
1	G	81	ASP
1	G	106	LEU
1	G	114	ARG
1	G	122	VAL
1	G	129	LEU
1	G	139	VAL
1	G	141	VAL
1	H	14	LEU
1	H	28	LEU
1	H	52	GLU
1	H	53	LEU
1	H	98	THR
1	H	100	LEU
1	H	106	LEU
1	H	114	ARG
1	H	122	VAL
1	H	139	VAL
1	I	14	LEU
1	I	16	THR
1	I	35	SER
1	I	92	MET
1	I	108	ARG

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Mol	Chain	Res	Type
1	I	114	ARG
1	I	117	LYS
1	I	122	VAL
1	J	16	THR
1	J	28	LEU
1	J	80	ASP
1	J	96	HIS
1	J	98	THR
1	J	114	ARG
1	J	129	LEU
1	K	16	THR
1	K	28	LEU
1	K	42	LEU
1	K	82	ILE
1	K	108	ARG
1	K	117	LYS
1	K	136	LYS
1	L	5	ILE
1	L	14	LEU
1	L	42	LEU
1	L	52	GLU
1	L	88	THR
1	L	102	THR
1	L	106	LEU
1	L	114	ARG
1	L	122	VAL
1	M	14	LEU
1	M	16	THR
1	M	18	MET
1	M	42	LEU
1	M	80	ASP
1	M	129	LEU
1	M	139	VAL
1	N	18	MET
1	N	85	ILE
1	N	92	MET
1	N	129	LEU
1	O	14	LEU
1	O	16	THR
1	O	27	VAL
1	O	28	LEU
1	O	82	ILE

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Mol	Chain	Res	Type
1	O	106	LEU
1	O	107	ARG
1	P	11	PRO
1	P	16	THR
1	P	28	LEU
1	P	42	LEU
1	P	102	THR
1	P	107	ARG
1	P	114	ARG
1	P	122	VAL
1	P	129	LEU
1	Q	27	VAL
1	Q	28	LEU
1	Q	42	LEU
1	Q	52	GLU
1	Q	82	ILE
1	Q	88	THR
1	Q	94	GLU
1	Q	98	THR
1	Q	100	LEU
1	Q	107	ARG
1	Q	122	VAL
1	Q	128	SER
1	Q	139	VAL
1	R	12	ASP
1	R	16	THR
1	R	42	LEU
1	R	76	VAL
1	R	81	ASP
1	R	122	VAL
1	R	128	SER
1	S	14	LEU
1	S	18	MET
1	S	35	SER
1	S	81	ASP
1	S	94	GLU
1	S	102	THR
1	S	106	LEU
1	S	107	ARG
1	S	114	ARG
1	S	122	VAL
1	S	141	VAL

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Mol	Chain	Res	Type
1	T	14	LEU
1	T	77	VAL
1	T	81	ASP
1	T	91	GLN
1	T	102	THR
1	T	114	ARG
1	T	122	VAL
1	T	129	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	49	ASN
1	D	75	ASN
1	D	96	HIS
1	F	96	HIS
1	G	49	ASN
1	G	91	GLN
1	I	72	GLN
1	K	49	ASN
1	K	72	GLN
1	L	49	ASN
1	O	72	GLN
1	P	4	HIS
1	Q	96	HIS
1	S	40	ASN
1	T	4	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 30 ligands modelled in this entry, 10 are monoatomic - leaving 20 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	FUL	J	1142	-	11,11,11	0.58	0	16,16,16	0.88	0
2	FUL	K	1142	-	11,11,11	0.52	0	16,16,16	0.99	1 (6%)
2	FUL	M	1142	-	11,11,11	0.59	0	16,16,16	0.99	0
2	FUL	H	1142	-	11,11,11	0.64	0	16,16,16	1.02	1 (6%)
2	FUL	Q	1142	-	11,11,11	0.59	0	16,16,16	0.90	0
2	FUL	P	1142	-	11,11,11	0.61	0	16,16,16	1.09	2 (12%)
2	FUL	T	1142	-	11,11,11	0.69	0	16,16,16	1.02	0
2	FUL	I	1144	-	11,11,11	0.58	0	16,16,16	1.13	1 (6%)
2	FUL	G	1142	-	11,11,11	0.56	0	16,16,16	0.65	0
2	FUL	A	1142	-	11,11,11	0.60	0	16,16,16	1.30	3 (18%)
2	FUL	R	1142	-	11,11,11	0.57	0	16,16,16	1.15	1 (6%)
2	FUL	C	1142	-	11,11,11	0.52	0	16,16,16	0.96	1 (6%)
2	FUL	L	1142	-	11,11,11	0.54	0	16,16,16	0.85	0
2	FUL	S	1142	-	11,11,11	0.58	0	16,16,16	1.05	1 (6%)
2	FUL	B	1142	-	11,11,11	0.65	0	16,16,16	0.93	1 (6%)
2	FUL	D	1142	-	11,11,11	0.58	0	16,16,16	0.81	0
2	FUL	F	1142	-	11,11,11	0.53	0	16,16,16	0.64	0
2	FUL	N	1142	-	11,11,11	0.57	0	16,16,16	0.91	0
2	FUL	O	1142	-	11,11,11	0.59	0	16,16,16	0.80	0
2	FUL	E	1141	-	11,11,11	0.61	0	16,16,16	1.14	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	FUL	J	1142	-	-	-	0/1/1/1
2	FUL	K	1142	-	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FUL	M	1142	-	-	-	0/1/1/1
2	FUL	H	1142	-	-	-	0/1/1/1
2	FUL	Q	1142	-	-	-	0/1/1/1
2	FUL	P	1142	-	-	-	0/1/1/1
2	FUL	T	1142	-	-	-	0/1/1/1
2	FUL	I	1144	-	-	-	0/1/1/1
2	FUL	G	1142	-	-	-	0/1/1/1
2	FUL	A	1142	-	-	-	0/1/1/1
2	FUL	R	1142	-	-	-	0/1/1/1
2	FUL	C	1142	-	-	-	0/1/1/1
2	FUL	L	1142	-	-	-	0/1/1/1
2	FUL	S	1142	-	-	-	0/1/1/1
2	FUL	B	1142	-	-	-	0/1/1/1
2	FUL	D	1142	-	-	-	0/1/1/1
2	FUL	F	1142	-	-	-	0/1/1/1
2	FUL	N	1142	-	-	-	0/1/1/1
2	FUL	O	1142	-	-	-	0/1/1/1
2	FUL	E	1141	-	-	-	0/1/1/1

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	S	1142	FUL	C1-O5-C5	-2.87	109.90	114.37
2	E	1141	FUL	C1-O5-C5	-2.72	110.14	114.37
2	A	1142	FUL	C1-O5-C5	-2.67	110.21	114.37
2	A	1142	FUL	O5-C5-C6	2.65	112.61	106.74
2	I	1144	FUL	O5-C5-C6	2.62	112.56	106.74
2	P	1142	FUL	C1-O5-C5	-2.60	110.33	114.37
2	B	1142	FUL	C1-O5-C5	-2.41	110.62	114.37
2	E	1141	FUL	O5-C5-C6	2.40	112.06	106.74
2	R	1142	FUL	O5-C5-C6	2.40	112.06	106.74
2	H	1142	FUL	C1-O5-C5	-2.34	110.74	114.37
2	K	1142	FUL	O5-C5-C6	2.33	111.91	106.74
2	P	1142	FUL	O5-C5-C6	2.31	111.87	106.74
2	A	1142	FUL	O5-C5-C4	-2.21	105.57	109.55
2	C	1142	FUL	C6-C5-C4	-2.15	109.14	113.08

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	T	1142	FUL	1	0
2	I	1144	FUL	2	0
2	C	1142	FUL	1	0
2	F	1142	FUL	3	0
2	E	1141	FUL	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	140/147 (95%)	-0.28	1 (0%) 84 81	26, 35, 46, 52	0
1	B	140/147 (95%)	-0.15	2 (1%) 73 70	27, 36, 54, 59	0
1	C	140/147 (95%)	-0.12	2 (1%) 73 70	28, 37, 57, 59	0
1	D	140/147 (95%)	-0.26	0 100 100	17, 35, 51, 55	1 (0%)
1	E	139/147 (94%)	-0.22	1 (0%) 84 81	24, 33, 49, 57	0
1	F	140/147 (95%)	0.02	1 (0%) 84 81	29, 41, 61, 64	0
1	G	140/147 (95%)	-0.06	0 100 100	29, 39, 56, 62	0
1	H	140/147 (95%)	-0.02	0 100 100	25, 40, 56, 58	0
1	I	142/147 (96%)	-0.24	0 100 100	28, 36, 51, 55	0
1	J	140/147 (95%)	0.07	1 (0%) 84 81	27, 38, 63, 68	0
1	K	140/147 (95%)	-0.20	2 (1%) 73 70	25, 36, 52, 56	0
1	L	140/147 (95%)	0.08	0 100 100	30, 43, 64, 66	0
1	M	140/147 (95%)	-0.20	0 100 100	26, 35, 50, 59	0
1	N	140/147 (95%)	-0.17	1 (0%) 84 81	24, 35, 46, 48	0
1	O	140/147 (95%)	-0.08	1 (0%) 84 81	27, 37, 58, 60	0
1	P	140/147 (95%)	-0.31	1 (0%) 84 81	25, 34, 44, 48	0
1	Q	140/147 (95%)	0.06	1 (0%) 84 81	30, 41, 66, 70	0
1	R	140/147 (95%)	-0.08	0 100 100	28, 39, 61, 62	0
1	S	140/147 (95%)	-0.29	0 100 100	23, 32, 49, 57	0
1	T	140/147 (95%)	-0.36	2 (1%) 73 70	22, 30, 47, 52	0
All	All	2801/2940 (95%)	-0.14	16 (0%) 85 83	17, 36, 57, 70	1 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	78	SER	3.7
1	T	79	GLY	3.3
1	O	78	SER	3.3
1	C	79	GLY	3.2
1	B	78	SER	2.8
1	T	80	ASP	2.7
1	E	79	GLY	2.7
1	K	79	GLY	2.7
1	K	141	VAL	2.7
1	F	79	GLY	2.6
1	N	78	SER	2.3
1	A	141	VAL	2.1
1	J	81	ASP	2.1
1	P	78	SER	2.1
1	Q	141	VAL	2.0
1	B	79	GLY	2.0

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

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

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($\text{\AA}^2$)	Q<0.9
2	FUL	M	1142	11/11	0.87	0.12	42,44,45,46	0
2	FUL	Q	1142	11/11	0.87	0.11	47,49,49,49	0
2	FUL	H	1142	11/11	0.88	0.11	33,36,38,38	0
2	FUL	J	1142	11/11	0.88	0.10	48,49,50,51	0
2	FUL	L	1142	11/11	0.88	0.11	41,44,45,46	0
2	FUL	A	1142	11/11	0.88	0.10	27,29,30,30	0
2	FUL	C	1142	11/11	0.88	0.10	35,38,39,39	0
2	FUL	I	1144	11/11	0.89	0.10	43,43,44,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FUL	F	1142	11/11	0.89	0.10	46,46,46,46	0
2	FUL	N	1142	11/11	0.89	0.09	33,35,35,35	0
2	FUL	K	1142	11/11	0.89	0.09	39,40,42,43	0
2	FUL	S	1142	11/11	0.89	0.11	33,35,37,38	0
2	FUL	R	1142	11/11	0.90	0.09	33,34,36,36	0
2	FUL	D	1142	11/11	0.91	0.10	37,37,38,40	0
2	FUL	G	1142	11/11	0.91	0.09	40,41,41,41	0
2	FUL	T	1142	11/11	0.91	0.09	37,37,38,40	0
2	FUL	O	1142	11/11	0.92	0.09	40,40,41,41	0
2	FUL	E	1141	11/11	0.93	0.08	33,34,35,35	0
2	FUL	B	1142	11/11	0.94	0.09	40,41,42,42	0
2	FUL	P	1142	11/11	0.94	0.06	36,38,39,39	0
3	K	C	1143	1/1	0.94	0.08	58,58,58,58	0
3	K	E	1142	1/1	0.94	0.04	56,56,56,56	0
3	K	Q	1143	1/1	0.95	0.05	44,44,44,44	0
3	K	B	1143	1/1	0.96	0.08	36,36,36,36	0
3	K	S	1143	1/1	0.96	0.04	59,59,59,59	0
3	K	L	1143	1/1	0.98	0.07	46,46,46,46	0
3	K	N	1143	1/1	0.98	0.02	47,47,47,47	0
3	K	J	1143	1/1	0.98	0.11	47,47,47,47	0
3	K	K	1143	1/1	0.98	0.04	47,47,47,47	0
3	K	D	1143	1/1	0.99	0.04	43,43,43,43	0

6.5 Other polymers

There are no such residues in this entry.