



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:41 PM UTC

PDB ID : 5FJW / pdb_00005fjw
Title : Yeast delta-COP-I mu-homology domain complexed with Dsl1 WxWx(MSE) peptide
Authors : Suckling, R.J.; Evans, P.R.; Owen, D.J.
Deposited on : 2015-10-13
Resolution : 2.80 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

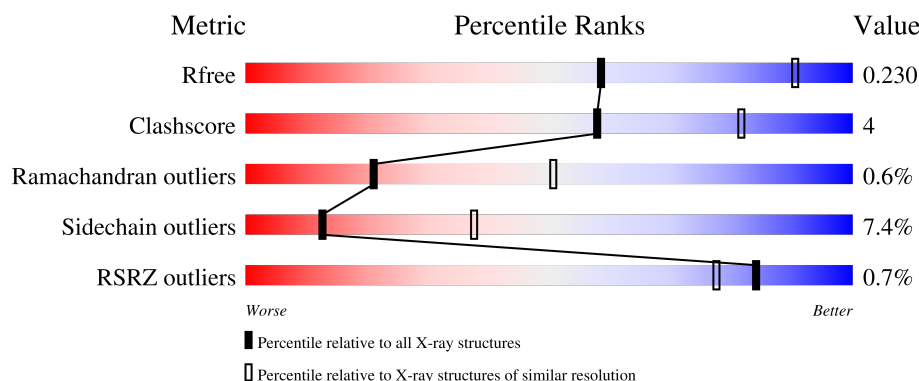
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

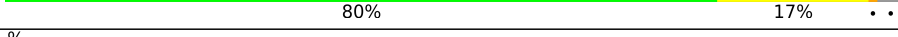
The reported resolution of this entry is 2.80 Å.

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.



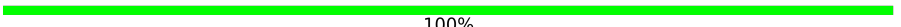
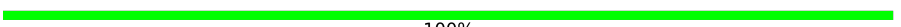
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	270	 81% 15% . .
1	B	270	 83% 13% . .
1	C	270	 80% 17% . .
1	D	270	 80% 17% . .
1	E	270	 81% 14% . .

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Mol	Chain	Length	Quality of chain
1	F	270	 81% 16% . .
1	G	270	 3% 79% 17% . .
1	H	270	 % 74% 21% . .
2	L	9	 100%
2	M	9	 89% 11%
2	N	9	 100%
2	O	9	 89% 11%
2	P	9	 78% 11% 11%
2	Q	9	 78% 22%
2	R	9	 78% 22%
2	S	9	 89% 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17707 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 COATOMER SUBUNIT DELTA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	Se	0	4	0
			2066	1299	337	423	1	6			
1	B	263	Total	C	N	O	S	Se	0	2	0
			2052	1289	333	423	1	6			
1	C	263	Total	C	N	O	S	Se	0	1	0
			2047	1286	332	422	1	6			
1	D	263	Total	C	N	O	S	Se	0	4	0
			2066	1299	334	426	1	6			
1	E	263	Total	C	N	O	S	Se	0	1	0
			2047	1286	332	422	1	6			
1	F	263	Total	C	N	O	S	Se	0	3	0
			2059	1294	335	423	1	6			
1	G	263	Total	C	N	O	S	Se	0	3	0
			2059	1294	334	424	1	6			
1	H	263	Total	C	N	O	S	Se	0	3	0
			2061	1296	336	422	1	6			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	277	GLY	-	expression tag	UNP P43621
A	278	PRO	-	expression tag	UNP P43621
A	279	LEU	-	expression tag	UNP P43621
A	280	GLY	-	expression tag	UNP P43621
A	281	SER	-	expression tag	UNP P43621
A	282	GLU	-	expression tag	UNP P43621
A	283	GLU	-	expression tag	UNP P43621
A	284	ASP	-	expression tag	UNP P43621
A	285	VAL	-	expression tag	UNP P43621
A	286	PRO	-	expression tag	UNP P43621
A	287	GLU	-	expression tag	UNP P43621
A	404	ALA	TRP	engineered mutation	UNP P43621
B	277	GLY	-	expression tag	UNP P43621

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Chain	Residue	Modelled	Actual	Comment	Reference
B	278	PRO	-	expression tag	UNP P43621
B	279	LEU	-	expression tag	UNP P43621
B	280	GLY	-	expression tag	UNP P43621
B	281	SER	-	expression tag	UNP P43621
B	282	GLU	-	expression tag	UNP P43621
B	283	GLU	-	expression tag	UNP P43621
B	284	ASP	-	expression tag	UNP P43621
B	285	VAL	-	expression tag	UNP P43621
B	286	PRO	-	expression tag	UNP P43621
B	287	GLU	-	expression tag	UNP P43621
B	404	ALA	TRP	engineered mutation	UNP P43621
C	277	GLY	-	expression tag	UNP P43621
C	278	PRO	-	expression tag	UNP P43621
C	279	LEU	-	expression tag	UNP P43621
C	280	GLY	-	expression tag	UNP P43621
C	281	SER	-	expression tag	UNP P43621
C	282	GLU	-	expression tag	UNP P43621
C	283	GLU	-	expression tag	UNP P43621
C	284	ASP	-	expression tag	UNP P43621
C	285	VAL	-	expression tag	UNP P43621
C	286	PRO	-	expression tag	UNP P43621
C	287	GLU	-	expression tag	UNP P43621
C	404	ALA	TRP	engineered mutation	UNP P43621
D	277	GLY	-	expression tag	UNP P43621
D	278	PRO	-	expression tag	UNP P43621
D	279	LEU	-	expression tag	UNP P43621
D	280	GLY	-	expression tag	UNP P43621
D	281	SER	-	expression tag	UNP P43621
D	282	GLU	-	expression tag	UNP P43621
D	283	GLU	-	expression tag	UNP P43621
D	284	ASP	-	expression tag	UNP P43621
D	285	VAL	-	expression tag	UNP P43621
D	286	PRO	-	expression tag	UNP P43621
D	287	GLU	-	expression tag	UNP P43621
D	404	ALA	TRP	engineered mutation	UNP P43621
E	277	GLY	-	expression tag	UNP P43621
E	278	PRO	-	expression tag	UNP P43621
E	279	LEU	-	expression tag	UNP P43621
E	280	GLY	-	expression tag	UNP P43621
E	281	SER	-	expression tag	UNP P43621
E	282	GLU	-	expression tag	UNP P43621
E	283	GLU	-	expression tag	UNP P43621

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Chain	Residue	Modelled	Actual	Comment	Reference
E	284	ASP	-	expression tag	UNP P43621
E	285	VAL	-	expression tag	UNP P43621
E	286	PRO	-	expression tag	UNP P43621
E	287	GLU	-	expression tag	UNP P43621
E	404	ALA	TRP	engineered mutation	UNP P43621
F	277	GLY	-	expression tag	UNP P43621
F	278	PRO	-	expression tag	UNP P43621
F	279	LEU	-	expression tag	UNP P43621
F	280	GLY	-	expression tag	UNP P43621
F	281	SER	-	expression tag	UNP P43621
F	282	GLU	-	expression tag	UNP P43621
F	283	GLU	-	expression tag	UNP P43621
F	284	ASP	-	expression tag	UNP P43621
F	285	VAL	-	expression tag	UNP P43621
F	286	PRO	-	expression tag	UNP P43621
F	287	GLU	-	expression tag	UNP P43621
F	404	ALA	TRP	engineered mutation	UNP P43621
G	277	GLY	-	expression tag	UNP P43621
G	278	PRO	-	expression tag	UNP P43621
G	279	LEU	-	expression tag	UNP P43621
G	280	GLY	-	expression tag	UNP P43621
G	281	SER	-	expression tag	UNP P43621
G	282	GLU	-	expression tag	UNP P43621
G	283	GLU	-	expression tag	UNP P43621
G	284	ASP	-	expression tag	UNP P43621
G	285	VAL	-	expression tag	UNP P43621
G	286	PRO	-	expression tag	UNP P43621
G	287	GLU	-	expression tag	UNP P43621
G	404	ALA	TRP	engineered mutation	UNP P43621
H	277	GLY	-	expression tag	UNP P43621
H	278	PRO	-	expression tag	UNP P43621
H	279	LEU	-	expression tag	UNP P43621
H	280	GLY	-	expression tag	UNP P43621
H	281	SER	-	expression tag	UNP P43621
H	282	GLU	-	expression tag	UNP P43621
H	283	GLU	-	expression tag	UNP P43621
H	284	ASP	-	expression tag	UNP P43621
H	285	VAL	-	expression tag	UNP P43621
H	286	PRO	-	expression tag	UNP P43621
H	287	GLU	-	expression tag	UNP P43621
H	404	ALA	TRP	engineered mutation	UNP P43621

- Molecule 2 is a protein called PROTEIN TRANSPORT PROTEIN DSL1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	9	Total	C	N	O	Se	0	0	0
			87	53	12	21	1			
2	M	9	Total	C	N	O	Se	0	0	0
			87	53	12	21	1			
2	N	9	Total	C	N	O	Se	0	0	0
			87	53	12	21	1			
2	O	9	Total	C	N	O	Se	0	0	0
			87	53	12	21	1			
2	P	9	Total	C	N	O	Se	0	0	0
			87	53	12	21	1			
2	Q	9	Total	C	N	O	Se	0	0	0
			87	53	12	21	1			
2	R	9	Total	C	N	O	Se	0	0	0
			87	53	12	21	1			
2	S	9	Total	C	N	O	Se	0	0	0
			87	53	12	21	1			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	417	MSE	VAL	conflict	UNP P53847
M	417	MSE	VAL	conflict	UNP P53847
N	417	MSE	VAL	conflict	UNP P53847
O	417	MSE	VAL	conflict	UNP P53847
P	417	MSE	VAL	conflict	UNP P53847
Q	417	MSE	VAL	conflict	UNP P53847
R	417	MSE	VAL	conflict	UNP P53847
S	417	MSE	VAL	conflict	UNP P53847

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	100	Total	O	0	0
			100	100		
3	B	106	Total	O	0	0
			106	106		
3	C	90	Total	O	0	0
			90	90		
3	D	97	Total	O	0	0
			97	97		
3	E	50	Total	O	0	0
			50	50		
3	F	42	Total	O	0	0
			42	42		

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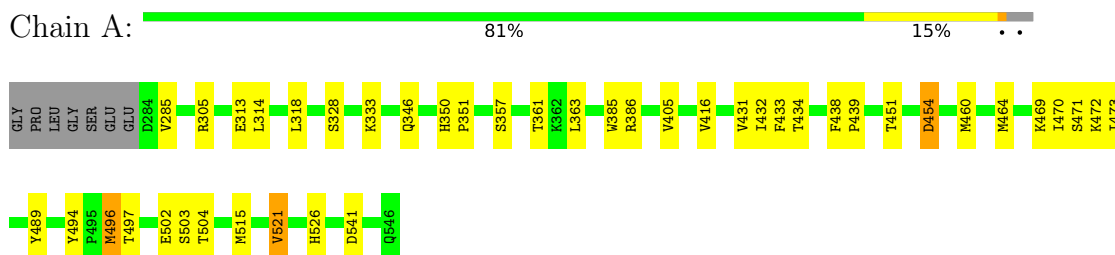
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	43	Total 43	O 43	0	0
3	H	21	Total 21	O 21	0	0
3	L	3	Total 3	O 3	0	0
3	M	2	Total 2	O 2	0	0

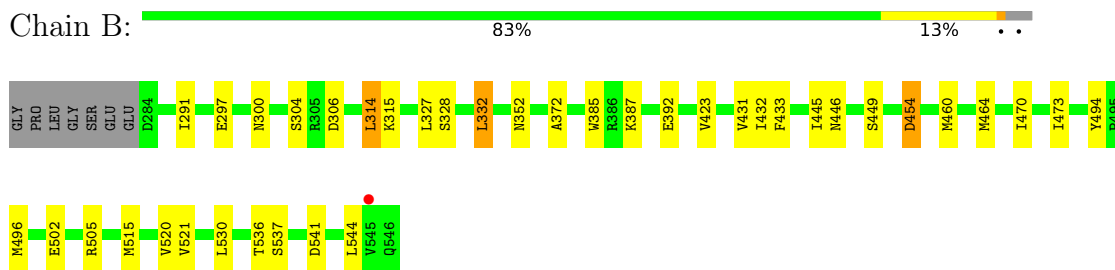
3 Residue-property plots

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.

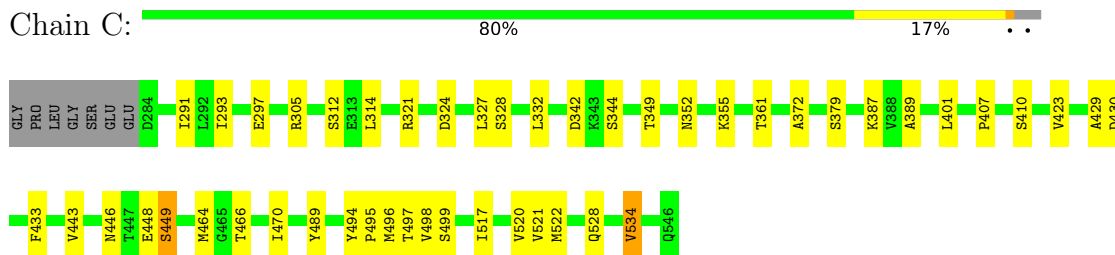
• Molecule 1: COATOMER SUBUNIT DELTA



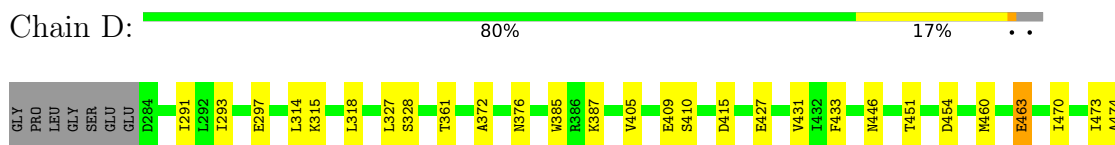
• Molecule 1: COATOMER SUBUNIT DELTA



• Molecule 1: COATOMER SUBUNIT DELTA

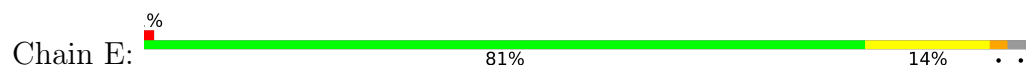


• Molecule 1: COATOMER SUBUNIT DELTA

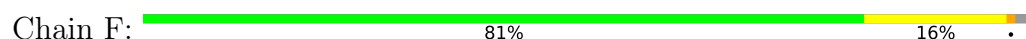




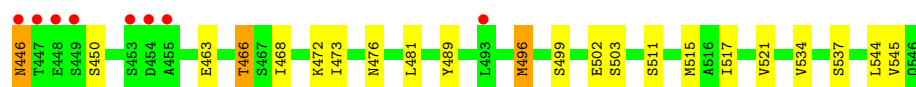
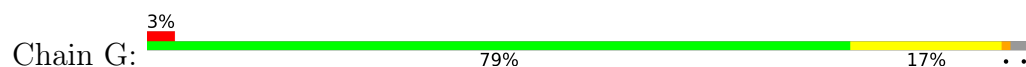
• Molecule 1: COATOMER SUBUNIT DELTA



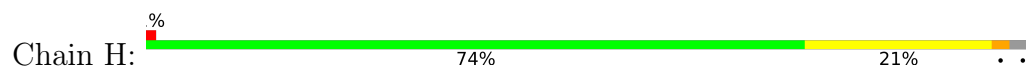
• Molecule 1: COATOMER SUBUNIT DELTA



• Molecule 1: COATOMER SUBUNIT DELTA



• Molecule 1: COATOMER SUBUNIT DELTA



• Molecule 2: PROTEIN TRANSPORT PROTEIN DSL1



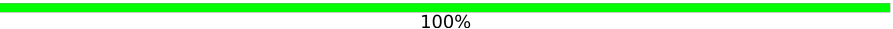
There are no outlier residues recorded for this chain.

- Molecule 2: PROTEIN TRANSPORT PROTEIN DSL1

Chain M:  89% 11%



- Molecule 2: PROTEIN TRANSPORT PROTEIN DSL1

Chain N:  100%

There are no outlier residues recorded for this chain.

- Molecule 2: PROTEIN TRANSPORT PROTEIN DSL1

Chain O:  89% 11%



- Molecule 2: PROTEIN TRANSPORT PROTEIN DSL1

Chain P:  78% 11% 11%



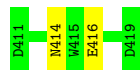
- Molecule 2: PROTEIN TRANSPORT PROTEIN DSL1

Chain Q:  78% 22%



- Molecule 2: PROTEIN TRANSPORT PROTEIN DSL1

Chain R:  78% 22%



- Molecule 2: PROTEIN TRANSPORT PROTEIN DSL1

Chain S:  89% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	68.41Å 79.30Å 163.35Å 79.71° 87.50° 83.60°	Depositor
Resolution (Å)	160.67 – 2.80 160.67 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.8 (160.67-2.80) 97.8 (160.67-2.80)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.82Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.170 , 0.239 (Not available) , 0.230	Depositor DCC
R_{free} test set	4059 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	73.8	Xtriage
Anisotropy	0.316	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 60.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17707	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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 i

5.1 Standard geometry i

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	0.94	1/2107 (0.0%)	1.02	2/2849 (0.1%)
1	B	0.93	0/2087	1.00	4/2824 (0.1%)
1	C	0.92	1/2079 (0.0%)	1.02	1/2813 (0.0%)
1	D	0.87	0/2107	1.02	2/2849 (0.1%)
1	E	0.80	1/2079 (0.0%)	0.97	1/2813 (0.0%)
1	F	0.81	1/2098 (0.0%)	0.94	1/2839 (0.0%)
1	G	0.81	0/2098	0.94	3/2839 (0.1%)
1	H	0.81	0/2101	0.95	5/2843 (0.2%)
2	L	0.92	0/89	1.06	0/118
2	M	0.72	0/89	0.86	0/118
2	N	0.99	0/89	0.83	0/118
2	O	0.83	0/89	1.13	0/118
2	P	0.74	0/89	1.02	0/118
2	Q	0.80	0/89	0.81	0/118
2	R	0.65	0/89	0.82	0/118
2	S	0.77	0/89	0.77	0/118
All	All	0.86	4/17468 (0.0%)	0.98	19/23613 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	D	0	2
1	H	0	1
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	541	ASP	CA-C	9.08	1.56	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	541	ASP	CA-C	6.13	1.55	1.52
1	A	541	ASP	CA-C	5.10	1.54	1.52
1	C	449	SER	CA-C	5.00	1.59	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	544	LEU	N-CA-C	7.37	120.56	108.99
1	E	544	LEU	N-CA-C	6.67	119.22	109.07
1	H	285	VAL	CA-C-N	6.41	126.43	119.89
1	H	285	VAL	C-N-CA	6.41	126.43	119.89
1	F	328	SER	N-CA-C	6.33	118.17	111.28
1	H	397	ILE	N-CA-C	6.31	113.80	107.55
1	A	496	MSE	N-CA-CB	6.25	119.78	110.29
1	D	410	SER	N-CA-C	6.03	118.82	111.82
1	A	464	MSE	N-CA-C	6.02	118.81	111.82
1	B	352	ASN	N-CA-C	5.61	118.92	111.75
1	G	410	SER	N-CA-C	5.58	120.14	113.23
1	B	544	LEU	N-CA-C	5.52	117.46	109.07
1	H	544	LEU	N-CA-C	5.44	117.12	108.79
1	H	505	ARG	N-CA-C	5.26	119.80	113.17
1	D	415	ASP	N-CA-C	-5.13	102.10	110.20
1	B	432	ILE	N-CA-C	5.08	115.30	107.77
1	C	410	SER	N-CA-C	5.07	117.54	111.71
1	G	352	ASN	N-CA-C	5.04	118.92	112.87
1	B	537	SER	N-CA-C	5.04	117.04	109.23

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	285	VAL	Peptide
1	D	446	ASN	Peptide
1	D	454	ASP	Peptide
1	H	285	VAL	Peptide

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	2066	0	2048	23	0
1	B	2052	0	2022	16	0
1	C	2047	0	2016	19	0
1	D	2066	0	2035	18	0
1	E	2047	0	2016	18	0
1	F	2059	0	2029	15	0
1	G	2059	0	2027	18	0
1	H	2061	0	2030	27	0
2	L	87	0	58	0	0
2	M	87	0	58	0	0
2	N	87	0	58	0	0
2	O	87	0	58	0	0
2	P	87	0	58	1	0
2	Q	87	0	58	0	0
2	R	87	0	58	0	0
2	S	87	0	58	0	0
3	A	100	0	0	1	0
3	B	106	0	0	0	0
3	C	90	0	0	1	0
3	D	97	0	0	1	0
3	E	50	0	0	2	0
3	F	42	0	0	0	0
3	G	43	0	0	1	0
3	H	21	0	0	0	0
3	L	3	0	0	0	0
3	M	2	0	0	0	0
All	All	17707	0	16687	152	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 (152) 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:H:468:ILE:HG21	1:H:481:LEU:HD21	1.74	0.69
1:A:494:TYR:CZ	1:A:496:MSE:HE3	2.32	0.65
1:A:494:TYR:CE1	1:A:496:MSE:HE3	2.32	0.65
1:H:403:THR:HG21	1:H:496:MSE:CE	2.26	0.64
1:A:346:GLN:HE22	1:A:386[A]:ARG:HE	1.45	0.63
1:B:327:LEU:CD2	1:B:530:LEU:HD22	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:445:ILE:HD11	1:H:457:VAL:HG23	1.83	0.61
1:G:468:ILE:HG21	1:G:481:LEU:HD21	1.83	0.60
1:F:468:ILE:HG21	1:F:481:LEU:HD21	1.83	0.60
1:C:293:ILE:HD11	1:C:332:LEU:HD22	1.82	0.60
1:B:328:SER:HB3	1:B:372:ALA:HB1	1.84	0.59
1:B:291:ILE:HD13	1:B:520:VAL:HG11	1.83	0.59
1:B:446[A]:ASN:CG	1:B:449:SER:HB3	2.28	0.58
1:F:297:GLU:HB3	1:F:314:LEU:HD11	1.84	0.58
1:C:328:SER:HB3	1:C:372:ALA:HB1	1.85	0.58
1:A:431:VAL:HG21	1:A:473:ILE:HD12	1.84	0.58
1:C:305:ARG:HG2	1:C:494:TYR:CE1	2.39	0.57
1:F:328:SER:HB3	1:F:372:ALA:HB1	1.88	0.56
1:H:438:PHE:CE1	1:H:460:MSE:HE2	2.41	0.55
1:A:333:LYS:HB2	3:A:2026:HOH:O	2.05	0.55
1:C:297:GLU:HB3	1:C:314:LEU:HD11	1.89	0.55
1:C:433:PHE:CE1	1:C:470:ILE:HD12	2.42	0.54
1:D:431:VAL:HG21	1:D:473:ILE:HD12	1.89	0.54
1:C:433:PHE:HE1	1:C:470:ILE:HD12	1.72	0.54
1:G:403:THR:HG21	1:G:496:MSE:SE	2.57	0.54
1:B:454:ASP:OD1	1:B:454:ASP:N	2.40	0.54
1:B:494:TYR:CZ	1:B:496:MSE:HE3	2.43	0.54
1:B:431:VAL:HG21	1:B:473:ILE:HD12	1.90	0.54
1:D:507:LYS:HA	1:D:534:VAL:HG13	1.90	0.53
1:G:295:ILE:HG21	1:G:511:SER:HB3	1.91	0.53
1:A:439:PRO:HD2	1:A:460:MSE:HE3	1.91	0.53
1:H:405:VAL:HG13	1:H:414:PHE:HB3	1.90	0.53
1:A:433:PHE:HE1	1:A:470:ILE:HD12	1.73	0.53
2:P:411:ASP:O	2:P:412:ASP:C	2.52	0.52
1:H:403:THR:HG21	1:H:496:MSE:HE1	1.92	0.52
1:G:435:ILE:HD12	1:G:496:MSE:HE2	1.91	0.51
1:A:405:VAL:HG22	1:A:416:VAL:HG22	1.91	0.51
1:E:376:ASN:ND2	1:E:378:GLN:HG3	2.25	0.51
1:G:399:LEU:HD11	1:G:420:TYR:HB2	1.92	0.51
1:C:293:ILE:O	1:C:534:VAL:HA	2.10	0.50
1:A:385:TRP:CZ2	1:A:515[A]:MSE:HG3	2.46	0.50
1:A:433:PHE:CE1	1:A:470:ILE:HD12	2.46	0.50
1:E:507:LYS:HA	1:E:534:VAL:HG13	1.92	0.49
1:G:431:VAL:HG21	1:G:473:ILE:HD12	1.95	0.49
1:D:297:GLU:HB3	1:D:314:LEU:HD11	1.93	0.49
1:E:328:SER:HB3	1:E:372:ALA:HB1	1.95	0.49
1:E:468:ILE:HG21	1:E:481:LEU:HD21	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:297:GLU:HB3	1:B:314:LEU:HD11	1.94	0.48
1:F:403:THR:CG2	1:F:496:MSE:HE3	2.43	0.48
1:H:328:SER:HB3	1:H:372:ALA:HB1	1.94	0.48
1:A:454:ASP:OD1	1:A:454:ASP:N	2.46	0.48
1:D:291:ILE:HD13	1:D:520:VAL:HG11	1.96	0.48
1:A:350:HIS:CE1	1:A:351:PRO:HD2	2.49	0.48
1:B:385:TRP:CZ2	1:B:515[A]:MSE:HG3	2.49	0.48
1:H:291:ILE:HD13	1:H:520:VAL:HG11	1.96	0.47
1:A:434:THR:CG2	1:A:497:THR:HB	2.44	0.47
1:E:435:ILE:HG23	1:E:496:MSE:HE2	1.96	0.47
1:F:443:VAL:HG21	1:F:466:THR:HG21	1.96	0.47
1:C:291:ILE:HD13	1:C:520:VAL:HG11	1.97	0.47
1:D:498:VAL:HG12	1:D:543:TYR:CD1	2.50	0.47
1:E:435:ILE:HD12	1:E:496:MSE:HE2	1.96	0.47
1:A:314:LEU:C	1:A:314:LEU:HD23	2.39	0.47
1:E:526:HIS:CD2	3:E:2047:HOH:O	2.68	0.46
1:F:390:PRO:HD3	1:H:390:PRO:HB3	1.97	0.46
1:G:443:VAL:HG21	1:G:466:THR:HG21	1.96	0.46
1:E:297:GLU:HB3	1:E:314:LEU:HD11	1.97	0.46
1:C:401:LEU:HD11	1:C:498:VAL:HG11	1.98	0.46
1:F:355:LYS:O	1:F:358:PHE:HB3	2.16	0.46
1:G:316:GLY:HA3	1:G:385:TRP:CE2	2.50	0.46
1:H:297:GLU:HB3	1:H:314:LEU:HD11	1.97	0.46
1:H:439:PRO:HD2	1:H:460:MSE:HE3	1.96	0.46
1:F:388:VAL:HG21	1:H:311:SER:CB	2.46	0.46
1:G:344:SER:O	1:G:387:LYS:HG3	2.15	0.46
1:H:329[B]:HIS:CE1	1:H:524:ASN:HD21	2.34	0.46
1:B:304:SER:HB2	1:B:306:ASP:OD1	2.16	0.45
1:B:327:LEU:HD22	1:B:530:LEU:HD22	1.95	0.45
1:C:293:ILE:HG21	1:C:517:ILE:HG13	1.97	0.45
1:C:443:VAL:HG21	1:C:466:THR:HG21	1.99	0.45
1:D:405:VAL:HG11	1:D:490:GLU:HG3	1.97	0.45
1:H:435:ILE:HD13	1:H:496:MSE:HG2	1.97	0.45
1:C:446:ASN:CG	1:C:449:SER:HB3	2.41	0.45
1:G:297:GLU:HB3	1:G:314:LEU:HD11	1.98	0.45
1:A:438:PHE:CD1	1:A:460:MSE:HE2	2.52	0.45
1:C:342:ASP:OD1	1:C:344:SER:OG	2.27	0.45
1:H:312:SER:OG	1:H:389:ALA:HB3	2.17	0.45
1:A:313:GLU:OE2	1:A:386[A]:ARG:NH1	2.44	0.44
1:A:432:ILE:HD11	1:A:469[B]:LYS:CE	2.48	0.44
1:D:497:THR:HG23	3:D:2088:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:515[B]:MSE:HE2	1:D:515[B]:MSE:HB2	1.87	0.44
1:E:399:LEU:HD11	1:E:420:TYR:HB2	1.97	0.44
1:E:332:LEU:HD23	1:E:332:LEU:N	2.33	0.44
1:D:474:ALA:HB3	1:D:477:ASP:HB3	2.00	0.44
1:H:435:ILE:HD13	1:H:496:MSE:CG	2.48	0.44
1:B:314:LEU:C	1:B:314:LEU:HD12	2.43	0.44
1:E:316:GLY:HA3	1:E:385:TRP:CE2	2.53	0.44
1:F:397:ILE:HG23	1:F:543:TYR:OH	2.18	0.44
1:G:515[B]:MSE:HE2	1:G:515[B]:MSE:HB2	1.93	0.44
1:H:420:TYR:CE1	1:H:479:GLY:HA3	2.53	0.44
1:A:432:ILE:HD11	1:A:469[B]:LYS:HE2	2.00	0.43
1:D:433:PHE:CE1	1:D:470:ILE:HD12	2.53	0.43
1:G:435:ILE:HD12	1:G:496:MSE:CE	2.48	0.43
1:G:468:ILE:CG2	1:G:481:LEU:HD21	2.47	0.43
1:C:494:TYR:CZ	1:C:496:MSE:HE3	2.53	0.43
1:H:336:ASP:O	1:H:362:LYS:NZ	2.52	0.43
1:E:329:HIS:NE2	1:E:372:ALA:HB2	2.34	0.43
1:H:515[B]:MSE:HB2	1:H:515[B]:MSE:HE2	1.66	0.43
1:B:300:ASN:OD1	1:B:541:ASP:HB3	2.19	0.43
1:D:385:TRP:CZ2	1:D:515[A]:MSE:HG3	2.54	0.43
1:H:403:THR:HG21	1:H:496:MSE:SE	2.68	0.43
1:G:325[A]:HIS:CE1	3:G:2018:HOH:O	2.71	0.43
1:A:305:ARG:HD3	1:A:494:TYR:CE2	2.53	0.43
1:B:433:PHE:HE1	1:B:470:ILE:HD12	1.84	0.43
1:E:376:ASN:CG	1:E:378:GLN:HG2	2.44	0.43
1:C:324:ASP:HB3	1:C:327:LEU:HD12	2.01	0.42
1:G:293:ILE:HG21	1:G:517:ILE:HD11	2.00	0.42
1:G:328:SER:HB3	1:G:372:ALA:HB1	2.02	0.42
1:C:495:PRO:HD3	3:C:2085:HOH:O	2.19	0.42
1:E:392:GLU:HA	3:E:2007:HOH:O	2.19	0.42
1:G:446:ASN:O	1:G:450:SER:OG	2.33	0.42
1:E:312:SER:OG	1:E:389:ALA:HB3	2.18	0.42
1:H:431:VAL:HG21	1:H:473:ILE:HD12	2.01	0.42
1:D:293:ILE:HG23	1:D:318:LEU:HD11	2.01	0.42
1:H:506:ASP:OD1	1:H:507:LYS:N	2.53	0.42
1:D:327:LEU:CD2	1:D:530:LEU:HD22	2.50	0.42
1:F:435:ILE:CD1	1:F:496:MSE:SE	3.18	0.42
1:E:446:ASN:HB3	1:E:449:SER:HB3	2.01	0.41
1:E:468:ILE:HG21	1:E:481:LEU:HD11	2.03	0.41
1:F:431:VAL:HG21	1:F:473:ILE:HD12	2.02	0.41
1:A:305:ARG:HG2	1:A:494:TYR:CE1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:439:PRO:HD2	1:E:460:MSE:HE3	2.02	0.41
1:H:329[B]:HIS:NE2	1:H:369:LYS:O	2.51	0.41
1:D:376:ASN:OD1	1:D:376:ASN:N	2.53	0.41
1:D:463[B]:GLU:HA	1:D:463[B]:GLU:OE1	2.20	0.41
1:D:328:SER:HB3	1:D:372:ALA:HB1	2.03	0.41
1:F:388:VAL:HG21	1:H:311:SER:HB2	2.02	0.41
1:B:332:LEU:N	1:B:332:LEU:HD23	2.36	0.41
1:C:429:ALA:O	1:C:430:ASP:C	2.64	0.41
1:G:350:HIS:O	1:G:351:PRO:C	2.64	0.41
1:H:494:TYR:N	1:H:494:TYR:CD1	2.88	0.41
1:A:350:HIS:CG	1:A:351:PRO:CD	3.04	0.41
1:D:460:MSE:HE2	1:D:460:MSE:HB3	2.02	0.41
1:H:293:ILE:HG23	1:H:318:LEU:HD11	2.02	0.41
1:D:475:ALA:O	1:D:477:ASP:N	2.54	0.41
1:C:321:ARG:HG3	1:C:379:SER:HA	2.03	0.40
1:F:291:ILE:HD13	1:F:520:VAL:HG11	2.04	0.40
1:F:403:THR:HG21	1:F:496:MSE:HE3	2.03	0.40
1:F:298:VAL:O	1:F:314:LEU:HD12	2.21	0.40
1:A:361:THR:HG22	1:A:363:LEU:HB2	2.04	0.40
1:B:433:PHE:CE1	1:B:470:ILE:HD12	2.56	0.40
1:C:312:SER:OG	1:C:389:ALA:HB3	2.22	0.40
1:A:521:VAL:HG13	1:A:526:HIS:HA	2.04	0.40
1:H:463:GLU:HB3	1:H:464:MSE:HE3	2.04	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	265/270 (98%)	253 (96%)	12 (4%)	0	100	100
1	B	263/270 (97%)	249 (95%)	14 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	262/270 (97%)	248 (95%)	13 (5%)	1 (0%)	30	60
1	D	265/270 (98%)	248 (94%)	15 (6%)	2 (1%)	16	44
1	E	262/270 (97%)	242 (92%)	20 (8%)	0	100	100
1	F	264/270 (98%)	247 (94%)	16 (6%)	1 (0%)	30	60
1	G	264/270 (98%)	248 (94%)	13 (5%)	3 (1%)	11	36
1	H	264/270 (98%)	233 (88%)	27 (10%)	4 (2%)	8	28
2	L	7/9 (78%)	7 (100%)	0	0	100	100
2	M	7/9 (78%)	7 (100%)	0	0	100	100
2	N	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
2	O	7/9 (78%)	7 (100%)	0	0	100	100
2	P	7/9 (78%)	5 (71%)	1 (14%)	1 (14%)	0	0
2	Q	7/9 (78%)	7 (100%)	0	0	100	100
2	R	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
2	S	7/9 (78%)	7 (100%)	0	0	100	100
All	All	2165/2232 (97%)	2020 (93%)	133 (6%)	12 (1%)	21	51

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	352	ASN
1	D	476	ASN
1	G	446	ASN
1	G	476	ASN
2	P	412	ASP
1	H	446	ASN
1	D	528	GLN
1	H	362	LYS
1	H	455	ALA
1	H	478	ALA
1	G	463	GLU
1	F	290	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	239/236 (101%)	227 (95%)	12 (5%)	22	54
1	B	237/236 (100%)	223 (94%)	14 (6%)	18	48
1	C	236/236 (100%)	221 (94%)	15 (6%)	16	44
1	D	239/236 (101%)	223 (93%)	16 (7%)	15	42
1	E	236/236 (100%)	215 (91%)	21 (9%)	9	29
1	F	238/236 (101%)	215 (90%)	23 (10%)	8	25
1	G	238/236 (101%)	220 (92%)	18 (8%)	12	36
1	H	238/236 (101%)	217 (91%)	21 (9%)	9	29
2	L	9/8 (112%)	9 (100%)	0	100	100
2	M	9/8 (112%)	8 (89%)	1 (11%)	6	20
2	N	9/8 (112%)	9 (100%)	0	100	100
2	O	9/8 (112%)	8 (89%)	1 (11%)	6	20
2	P	9/8 (112%)	9 (100%)	0	100	100
2	Q	9/8 (112%)	7 (78%)	2 (22%)	1	3
2	R	9/8 (112%)	7 (78%)	2 (22%)	1	3
2	S	9/8 (112%)	8 (89%)	1 (11%)	6	20
All	All	1973/1952 (101%)	1826 (92%)	147 (8%)	13	37

All (147) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	318	LEU
1	A	328	SER
1	A	357	SER
1	A	451	THR
1	A	454	ASP
1	A	471	SER
1	A	472	LYS
1	A	489	TYR
1	A	502	GLU
1	A	503	SER
1	A	504	THR
1	A	521	VAL
1	B	314	LEU
1	B	315	LYS

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Mol	Chain	Res	Type
1	B	332	LEU
1	B	387	LYS
1	B	392	GLU
1	B	423	VAL
1	B	445	ILE
1	B	454	ASP
1	B	460	MSE
1	B	464	MSE
1	B	502	GLU
1	B	505	ARG
1	B	521	VAL
1	B	536	THR
1	C	349	THR
1	C	355	LYS
1	C	361	THR
1	C	387	LYS
1	C	407	PRO
1	C	423	VAL
1	C	448	GLU
1	C	464	MSE
1	C	489	TYR
1	C	497	THR
1	C	499	SER
1	C	521	VAL
1	C	522	MSE
1	C	528	GLN
1	C	534	VAL
1	D	315	LYS
1	D	361	THR
1	D	387	LYS
1	D	409[A]	GLU
1	D	409[B]	GLU
1	D	427	GLU
1	D	451	THR
1	D	463[A]	GLU
1	D	463[B]	GLU
1	D	502	GLU
1	D	503	SER
1	D	504	THR
1	D	505	ARG
1	D	521	VAL
1	D	529	GLU

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Mol	Chain	Res	Type
1	D	534	VAL
1	E	308	THR
1	E	315	LYS
1	E	328	SER
1	E	332	LEU
1	E	360	SER
1	E	361	THR
1	E	379	SER
1	E	402	THR
1	E	450	SER
1	E	464	MSE
1	E	489	TYR
1	E	496	MSE
1	E	501	GLN
1	E	502	GLU
1	E	503	SER
1	E	504	THR
1	E	521	VAL
1	E	534	VAL
1	E	537	SER
1	E	544	LEU
1	E	545	VAL
1	F	308	THR
1	F	315	LYS
1	F	332	LEU
1	F	333	LYS
1	F	355	LYS
1	F	360	SER
1	F	361	THR
1	F	406	SER
1	F	409	GLU
1	F	447	THR
1	F	448	GLU
1	F	452	CYS
1	F	486	GLU
1	F	496	MSE
1	F	497	THR
1	F	499	SER
1	F	502	GLU
1	F	504	THR
1	F	505	ARG
1	F	521	VAL

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Mol	Chain	Res	Type
1	F	533	ASP
1	F	534	VAL
1	F	536	THR
1	G	315	LYS
1	G	332	LEU
1	G	361	THR
1	G	363	LEU
1	G	388	VAL
1	G	406	SER
1	G	427	GLU
1	G	466	THR
1	G	472	LYS
1	G	489	TYR
1	G	496	MSE
1	G	499	SER
1	G	502	GLU
1	G	503	SER
1	G	521	VAL
1	G	534	VAL
1	G	537	SER
1	G	545	VAL
1	H	332	LEU
1	H	337	SER
1	H	361	THR
1	H	387	LYS
1	H	402	THR
1	H	405	VAL
1	H	409	GLU
1	H	426	THR
1	H	447	THR
1	H	448	GLU
1	H	452	CYS
1	H	469	LYS
1	H	489	TYR
1	H	496	MSE
1	H	499	SER
1	H	505	ARG
1	H	521	VAL
1	H	524	ASN
1	H	534	VAL
1	H	536	THR
1	H	537	SER

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Mol	Chain	Res	Type
2	M	418	GLU
2	O	416	GLU
2	Q	418	GLU
2	Q	419	ASP
2	R	414	ASN
2	R	416	GLU
2	S	417	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	331	ASN
1	A	346	GLN
1	A	518	GLN
1	A	528	GLN
1	A	546	GLN
1	B	459	ASN
1	B	524	ASN
1	C	289	ASN
1	C	329	HIS
1	C	331	ASN
1	C	411	GLN
1	C	462	GLN
1	C	476	ASN
1	C	546	GLN
1	D	528	GLN
1	E	331	ASN
1	E	446	ASN
1	E	459	ASN
1	E	476	ASN
1	F	310	HIS
1	F	331	ASN
1	F	346	GLN
1	F	476	ASN
1	F	524	ASN
1	G	329	HIS
1	G	331	ASN
1	H	288	ASN
1	H	346	GLN
1	H	462	GLN
1	H	476	ASN
1	H	524	ASN

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Mol	Chain	Res	Type
1	H	528	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	258/270 (95%)	-0.69	0 100 100	31, 62, 103, 142	3 (1%)
1	B	258/270 (95%)	-0.59	1 (0%) 88 84	36, 65, 106, 154	2 (0%)
1	C	258/270 (95%)	-0.44	0 100 100	26, 73, 122, 153	0
1	D	258/270 (95%)	-0.51	0 100 100	43, 72, 120, 161	3 (1%)
1	E	258/270 (95%)	-0.20	3 (1%) 76 68	53, 87, 170, 198	0
1	F	258/270 (95%)	-0.22	0 100 100	41, 92, 168, 188	2 (0%)
1	G	258/270 (95%)	-0.06	8 (3%) 51 41	46, 94, 177, 211	2 (0%)
1	H	258/270 (95%)	0.02	2 (0%) 82 75	57, 109, 184, 208	2 (0%)
2	L	8/9 (88%)	-0.24	0 100 100	59, 88, 120, 124	0
2	M	8/9 (88%)	-0.32	0 100 100	63, 95, 119, 121	0
2	N	8/9 (88%)	-0.06	0 100 100	76, 109, 130, 147	0
2	O	8/9 (88%)	-0.08	0 100 100	59, 103, 127, 132	0
2	P	8/9 (88%)	-0.62	0 100 100	83, 106, 112, 113	0
2	Q	8/9 (88%)	-0.20	0 100 100	73, 104, 118, 128	0
2	R	8/9 (88%)	-0.29	0 100 100	78, 106, 124, 137	0
2	S	8/9 (88%)	-0.11	0 100 100	104, 128, 136, 158	0
All	All	2128/2232 (95%)	-0.33	14 (0%) 84 77	26, 81, 157, 211	14 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	545	VAL	3.8
1	G	449	SER	3.6
1	G	453	SER	3.4
1	G	455	ALA	3.1
1	E	453	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	G	448	GLU	2.4
1	E	284	ASP	2.4
1	G	446	ASN	2.4
1	G	454	ASP	2.4
1	H	433	PHE	2.3
1	E	454	ASP	2.2
1	H	455	ALA	2.2
1	G	447	THR	2.2
1	G	493	LEU	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.