



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 7, 2026 – 01:10 AM UTC

PDB ID : 5OAT / pdb_00005oat
Title : PINK1 structure
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Deposited on : 2017-06-23
Resolution : 2.78 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

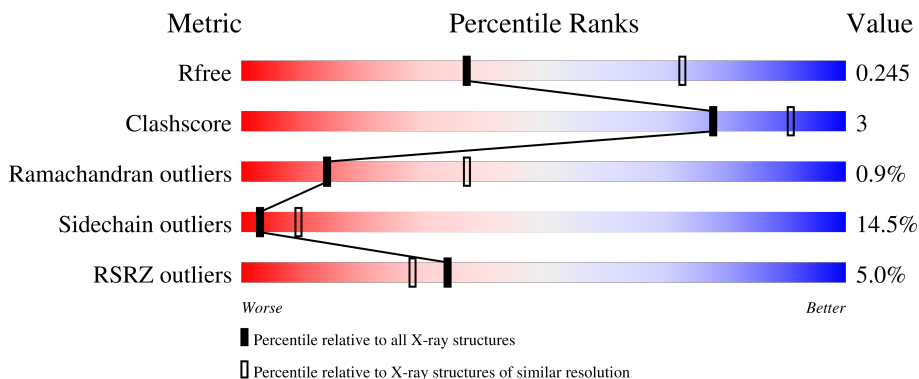
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.78 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.



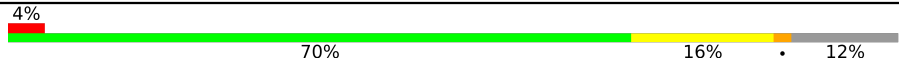
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	411	 3% 70% 16% 12%
1	B	411	 5% 71% 15% 12%
1	C	411	 5% 69% 16% 12%
1	D	411	 4% 70% 14% 12%
1	E	411	 6% 70% 15% 12%

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Mol	Chain	Length	Quality of chain
1	F	411	 <p>4% 70% 16% • 12%</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16758 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 Serine/threonine-protein kinase PINK1, mitochondrial-like Protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	360	2787	1793	464	512	1	17	0	0	0
1	B	360	2787	1793	464	512	1	17	0	0	0
1	C	360	2787	1793	464	512	1	17	0	0	0
1	D	360	2787	1793	464	512	1	17	0	0	0
1	E	360	2787	1793	464	512	1	17	0	0	0
1	F	360	2787	1793	464	512	1	17	0	0	0

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	205	GLU	SER	conflict	UNP D6WMX4
A	?	-	ILE	deletion	UNP D6WMX4
A	?	-	GLN	deletion	UNP D6WMX4
A	?	-	GLU	deletion	UNP D6WMX4
A	?	-	LEU	deletion	UNP D6WMX4
A	?	-	GLU	deletion	UNP D6WMX4
A	?	-	GLY	deletion	UNP D6WMX4
A	?	-	SER	deletion	UNP D6WMX4
A	?	-	LYS	deletion	UNP D6WMX4
A	?	-	ASP	deletion	UNP D6WMX4
A	?	-	LEU	deletion	UNP D6WMX4
A	527	ALA	GLU	conflict	UNP D6WMX4
A	528	ALA	LYS	conflict	UNP D6WMX4
B	205	GLU	SER	conflict	UNP D6WMX4
B	?	-	ILE	deletion	UNP D6WMX4
B	?	-	GLN	deletion	UNP D6WMX4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	GLU	deletion	UNP D6WMX4
B	?	-	LEU	deletion	UNP D6WMX4
B	?	-	GLU	deletion	UNP D6WMX4
B	?	-	GLY	deletion	UNP D6WMX4
B	?	-	SER	deletion	UNP D6WMX4
B	?	-	LYS	deletion	UNP D6WMX4
B	?	-	ASP	deletion	UNP D6WMX4
B	?	-	LEU	deletion	UNP D6WMX4
B	527	ALA	GLU	conflict	UNP D6WMX4
B	528	ALA	LYS	conflict	UNP D6WMX4
C	205	GLU	SER	conflict	UNP D6WMX4
C	?	-	ILE	deletion	UNP D6WMX4
C	?	-	GLN	deletion	UNP D6WMX4
C	?	-	GLU	deletion	UNP D6WMX4
C	?	-	LEU	deletion	UNP D6WMX4
C	?	-	GLU	deletion	UNP D6WMX4
C	?	-	GLY	deletion	UNP D6WMX4
C	?	-	SER	deletion	UNP D6WMX4
C	?	-	LYS	deletion	UNP D6WMX4
C	?	-	ASP	deletion	UNP D6WMX4
C	?	-	LEU	deletion	UNP D6WMX4
C	527	ALA	GLU	conflict	UNP D6WMX4
C	528	ALA	LYS	conflict	UNP D6WMX4
D	205	GLU	SER	conflict	UNP D6WMX4
D	?	-	ILE	deletion	UNP D6WMX4
D	?	-	GLN	deletion	UNP D6WMX4
D	?	-	GLU	deletion	UNP D6WMX4
D	?	-	LEU	deletion	UNP D6WMX4
D	?	-	GLU	deletion	UNP D6WMX4
D	?	-	GLY	deletion	UNP D6WMX4
D	?	-	SER	deletion	UNP D6WMX4
D	?	-	LYS	deletion	UNP D6WMX4
D	?	-	ASP	deletion	UNP D6WMX4
D	?	-	LEU	deletion	UNP D6WMX4
D	527	ALA	GLU	conflict	UNP D6WMX4
D	528	ALA	LYS	conflict	UNP D6WMX4
E	205	GLU	SER	conflict	UNP D6WMX4
E	?	-	ILE	deletion	UNP D6WMX4
E	?	-	GLN	deletion	UNP D6WMX4
E	?	-	GLU	deletion	UNP D6WMX4
E	?	-	LEU	deletion	UNP D6WMX4
E	?	-	GLU	deletion	UNP D6WMX4

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Chain	Residue	Modelled	Actual	Comment	Reference
E	?	-	GLY	deletion	UNP D6WMX4
E	?	-	SER	deletion	UNP D6WMX4
E	?	-	LYS	deletion	UNP D6WMX4
E	?	-	ASP	deletion	UNP D6WMX4
E	?	-	LEU	deletion	UNP D6WMX4
E	527	ALA	GLU	conflict	UNP D6WMX4
E	528	ALA	LYS	conflict	UNP D6WMX4
F	205	GLU	SER	conflict	UNP D6WMX4
F	?	-	ILE	deletion	UNP D6WMX4
F	?	-	GLN	deletion	UNP D6WMX4
F	?	-	GLU	deletion	UNP D6WMX4
F	?	-	LEU	deletion	UNP D6WMX4
F	?	-	GLU	deletion	UNP D6WMX4
F	?	-	GLY	deletion	UNP D6WMX4
F	?	-	SER	deletion	UNP D6WMX4
F	?	-	LYS	deletion	UNP D6WMX4
F	?	-	ASP	deletion	UNP D6WMX4
F	?	-	LEU	deletion	UNP D6WMX4
F	527	ALA	GLU	conflict	UNP D6WMX4
F	528	ALA	LYS	conflict	UNP D6WMX4

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	C	2	Total Mg 2 2	0	0
2	E	2	Total Mg 2 2	0	0
2	F	1	Total Mg 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	6	Total O 6 6	0	0
3	B	9	Total O 9 9	0	0
3	C	3	Total O 3 3	0	0

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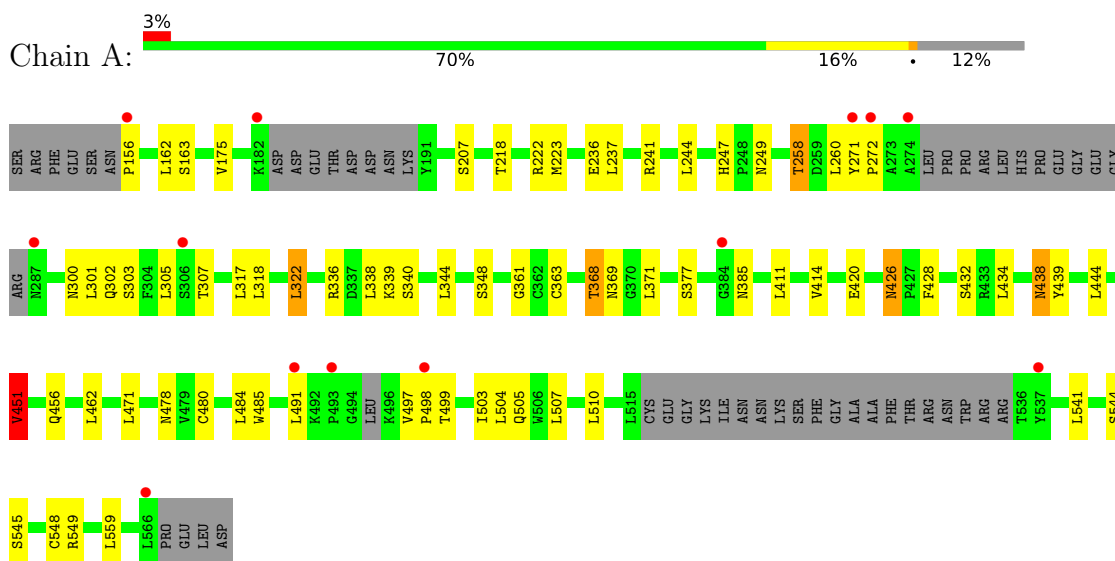
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	6	Total O 6 6	0	0
3	E	1	Total O 1 1	0	0
3	F	5	Total O 5 5	0	0

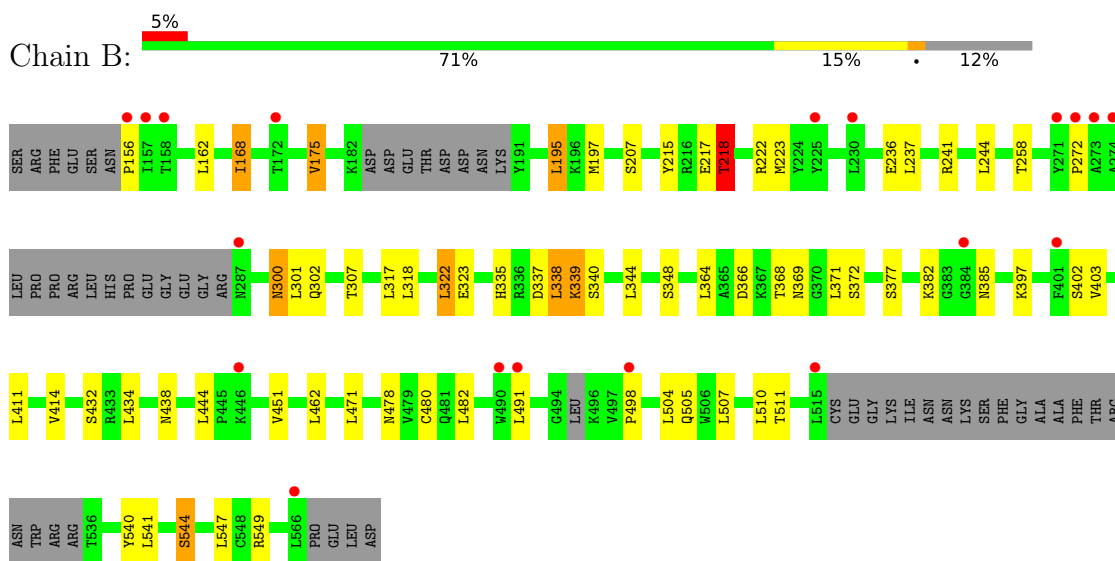
3 Residue-property plots [i](#)

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.

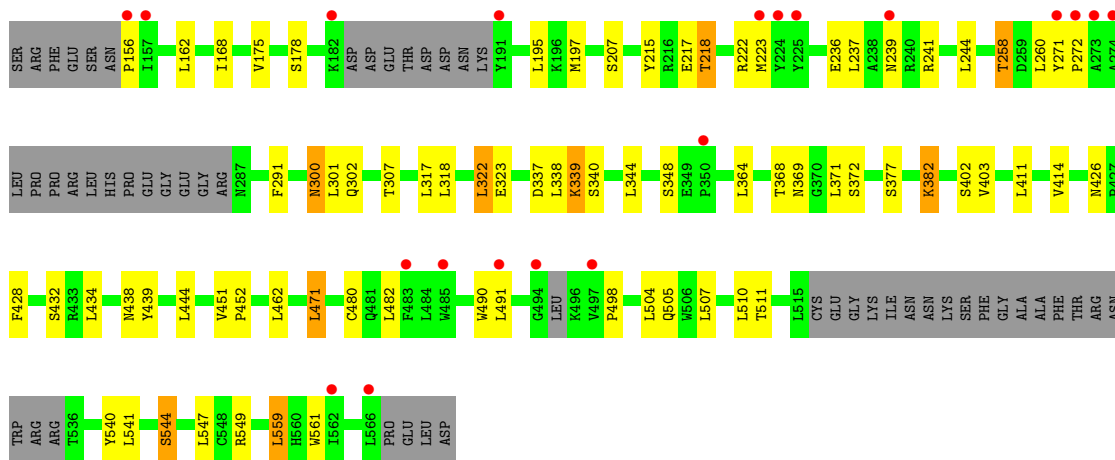
- Molecule 1: Serine/threonine-protein kinase PINK1, mitochondrial-like Protein



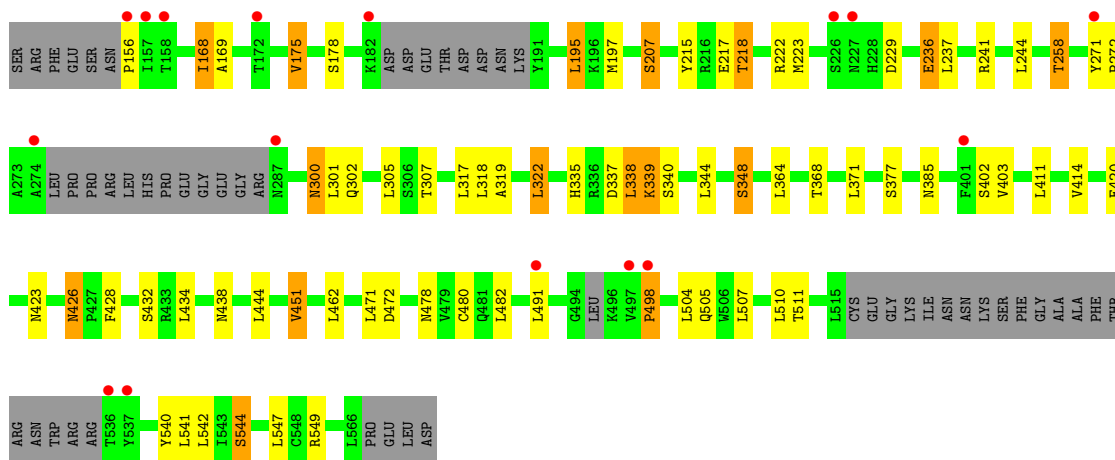
- Molecule 1: Serine/threonine-protein kinase PINK1, mitochondrial-like Protein



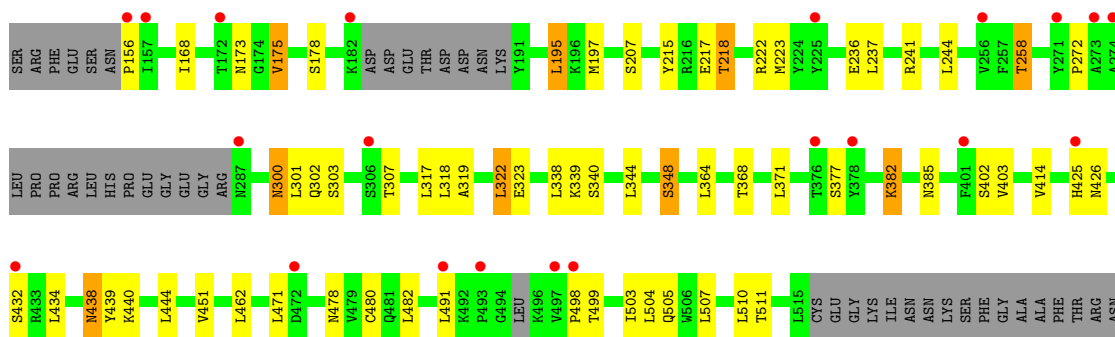
- Molecule 1: Serine/threonine-protein kinase PINK1, mitochondrial-like Protein



• Molecule 1: Serine/threonine-protein kinase PINK1, mitochondrial-like Protein

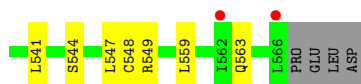
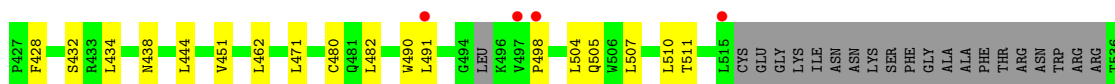
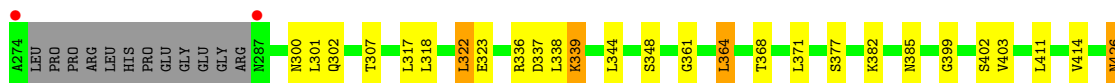
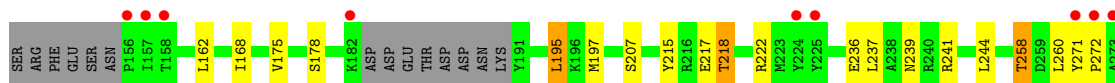


• Molecule 1: Serine/threonine-protein kinase PINK1, mitochondrial-like Protein





- Molecule 1: Serine/threonine-protein kinase PINK1, mitochondrial-like Protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.92Å 116.74Å 179.34Å 90.00° 94.29° 90.00°	Depositor
Resolution (Å)	178.84 – 2.78 178.84 – 2.78	Depositor EDS
% Data completeness (in resolution range)	99.8 (178.84-2.78) 99.8 (178.84-2.78)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.77Å)	Xtrriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.205 , 0.246 0.206 , 0.245	Depositor DCC
R_{free} test set	4195 reflections (4.14%)	wwPDB-VP
Wilson B-factor (Å ²)	73.7	Xtrriage
Anisotropy	0.192	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 42.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16758	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.87 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.7648e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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

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	0.96	1/2842 (0.0%)	1.13	10/3876 (0.3%)
1	B	1.03	0/2842	1.12	6/3876 (0.2%)
1	C	0.97	1/2842 (0.0%)	1.10	2/3876 (0.1%)
1	D	1.01	2/2842 (0.1%)	1.12	6/3876 (0.2%)
1	E	0.95	1/2842 (0.0%)	1.06	6/3876 (0.2%)
1	F	0.99	2/2842 (0.1%)	1.12	4/3876 (0.1%)
All	All	0.99	7/17052 (0.0%)	1.11	34/23256 (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	B	0	1
1	C	0	1
1	D	0	1
1	E	0	1
1	F	0	1
All	All	0	6

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	382	LYS	CA-C	7.06	1.61	1.52
1	F	399	GLY	C-O	-6.40	1.15	1.24
1	A	497	VAL	N-CA	5.99	1.50	1.46
1	E	382	LYS	CA-C	5.58	1.59	1.52
1	D	271	TYR	CA-C	5.50	1.59	1.52

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	497	VAL	N-CA-C	7.26	114.16	107.56
1	A	438	ASN	CB-CA-C	-6.64	97.56	110.11
1	D	438	ASN	CB-CA-C	-6.47	98.86	110.70
1	F	361	GLY	N-CA-C	6.32	121.42	114.40
1	D	175	VAL	CB-CA-C	-6.17	99.67	110.71

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	432	SER	Peptide
1	B	432	SER	Peptide
1	C	432	SER	Peptide
1	D	432	SER	Peptide
1	E	432	SER	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	2787	0	2741	15	0
1	B	2787	0	2741	14	0
1	C	2787	0	2741	18	0
1	D	2787	0	2741	22	0
1	E	2787	0	2741	17	0
1	F	2787	0	2741	14	0
2	A	1	0	0	0	0
2	C	2	0	0	0	0
2	E	2	0	0	0	0
2	F	1	0	0	0	0
3	A	6	0	0	0	0
3	B	9	0	0	0	0
3	C	3	0	0	0	0
3	D	6	0	0	2	0
3	E	1	0	0	0	0
3	F	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	16758	0	16446	93	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 3.

The worst 5 of 93 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:D:156:PRO:HG3	1:D:223:MET:HE2	1.79	0.65
1:A:260:LEU:C	1:A:271:TYR:N	2.55	0.65
1:A:322:LEU:HD13	1:A:414:VAL:HG11	1.80	0.64
1:B:168:ILE:O	1:C:239:ASN:ND2	2.31	0.64
1:F:322:LEU:HD13	1:F:414:VAL:HG11	1.81	0.63

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	347/411 (84%)	325 (94%)	20 (6%)	2 (1%)	21 47
1	B	347/411 (84%)	326 (94%)	18 (5%)	3 (1%)	14 37
1	C	347/411 (84%)	328 (94%)	15 (4%)	4 (1%)	10 30
1	D	347/411 (84%)	323 (93%)	20 (6%)	4 (1%)	10 30
1	E	347/411 (84%)	327 (94%)	17 (5%)	3 (1%)	14 37
1	F	347/411 (84%)	325 (94%)	19 (6%)	3 (1%)	14 37
All	All	2082/2466 (84%)	1954 (94%)	109 (5%)	19 (1%)	14 37

5 of 19 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	272	PRO
1	B	272	PRO
1	E	272	PRO
1	F	272	PRO
1	B	402	SER

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	300/360 (83%)	259 (86%)	41 (14%)	3 11
1	B	300/360 (83%)	255 (85%)	45 (15%)	3 9
1	C	300/360 (83%)	256 (85%)	44 (15%)	3 9
1	D	300/360 (83%)	257 (86%)	43 (14%)	3 10
1	E	300/360 (83%)	256 (85%)	44 (15%)	3 9
1	F	300/360 (83%)	256 (85%)	44 (15%)	3 9
All	All	1800/2160 (83%)	1539 (86%)	261 (14%)	3 9

5 of 261 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	317	LEU
1	F	364	LEU
1	F	548	CYS
1	C	301	LEU
1	C	244	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 64 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	302	GLN
1	F	405	ASN
1	C	243	HIS
1	C	239	ASN

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Mol	Chain	Res	Type
1	F	426	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

6 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	SEP	D	207	1	8,9,10	0.89	0	7,12,14	2.41	1 (14%)
1	SEP	F	207	1	8,9,10	0.78	0	7,12,14	2.55	1 (14%)
1	SEP	A	207	1	8,9,10	0.75	0	7,12,14	2.82	1 (14%)
1	SEP	E	207	1	8,9,10	0.72	0	7,12,14	2.46	3 (42%)
1	SEP	C	207	1	8,9,10	0.84	0	7,12,14	2.73	1 (14%)
1	SEP	B	207	1	8,9,10	0.84	0	7,12,14	2.69	3 (42%)

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	SEP	D	207	1	-	2/6/8/10	-
1	SEP	F	207	1	-	3/6/8/10	-
1	SEP	A	207	1	-	2/6/8/10	-
1	SEP	E	207	1	-	4/6/8/10	-
1	SEP	C	207	1	-	3/6/8/10	-
1	SEP	B	207	1	-	5/6/8/10	-

There are no bond length outliers.

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	207	SEP	OG-CB-CA	6.83	114.79	108.14
1	C	207	SEP	OG-CB-CA	6.70	114.66	108.14
1	F	207	SEP	OG-CB-CA	6.05	114.04	108.14
1	B	207	SEP	OG-CB-CA	5.98	113.96	108.14
1	D	207	SEP	OG-CB-CA	5.52	113.52	108.14

There are no chirality outliers.

5 of 19 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	207	SEP	N-CA-CB-OG
1	A	207	SEP	C-CA-CB-OG
1	B	207	SEP	N-CA-CB-OG
1	B	207	SEP	C-CA-CB-OG
1	B	207	SEP	CB-OG-P-O1P

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	207	SEP	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1
1	C	1
1	E	1
1	F	1
1	D	1
1	A	1

The worst 5 of 6 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	260:LEU	C	271:TYR	N	3.02
1	C	260:LEU	C	271:TYR	N	3.00
1	E	260:LEU	C	271:TYR	N	2.93
1	F	260:LEU	C	271:TYR	N	2.86
1	D	260:LEU	C	271:TYR	N	2.82

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	359/411 (87%)	0.07	13 (3%) 46 39	44, 76, 124, 157	0
1	B	359/411 (87%)	0.19	19 (5%) 32 26	43, 72, 119, 174	0
1	C	359/411 (87%)	0.28	20 (5%) 30 25	47, 81, 133, 189	0
1	D	359/411 (87%)	0.21	16 (4%) 38 32	42, 74, 124, 178	0
1	E	359/411 (87%)	0.34	23 (6%) 25 20	45, 80, 136, 166	0
1	F	359/411 (87%)	0.21	17 (4%) 36 31	44, 75, 123, 194	0
All	All	2154/2466 (87%)	0.21	108 (5%) 34 28	42, 76, 128, 194	0

The worst 5 of 108 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	172	THR	5.0
1	C	156	PRO	5.0
1	E	274	ALA	5.0
1	D	401	PHE	4.6
1	D	156	PRO	4.5

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	SEP	F	207	10/11	0.84	0.12	71,90,141,160	0
1	SEP	C	207	10/11	0.86	0.12	69,88,143,148	0
1	SEP	A	207	10/11	0.89	0.09	63,86,131,135	0
1	SEP	D	207	10/11	0.90	0.12	64,80,123,130	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
1	SEP	E	207	10/11	0.93	0.09	72,88,120,121	0
1	SEP	B	207	10/11	0.94	0.10	57,78,116,121	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(\AA^2)	Q<0.9
2	MG	E	602	1/1	0.86	0.33	98,98,98,98	0
2	MG	A	601	1/1	0.87	0.24	73,73,73,73	0
2	MG	C	602	1/1	0.90	0.30	92,92,92,92	0
2	MG	F	601	1/1	0.90	0.19	74,74,74,74	0
2	MG	C	601	1/1	0.92	0.24	68,68,68,68	0
2	MG	E	601	1/1	0.96	0.26	81,81,81,81	0

6.5 Other polymers [i](#)

There are no such residues in this entry.