



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:20 PM UTC

PDB ID : 5XJX / pdb_00005xjx
Title : Pre-formed plant receptor ERL1-TMM complex
Authors : Chai, J.; Lin, G.; Zhang, L.; Han, Z.; Yang, X.; Liu, W.; Qi, Y.; Chang, J.;
Li, E.
Deposited on : 2017-05-04
Resolution : 3.06 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
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

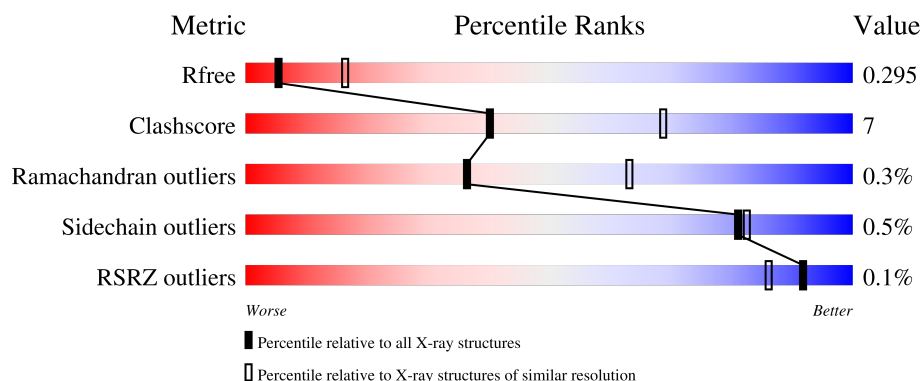
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.06 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	C	433	 70% 14% 16%
1	D	433	 69% 15% 16%
1	F	433	 68% 16% 16%
1	H	433	 69% 15% 16%
1	J	433	 72% 12% 16%

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Mol	Chain	Length	Quality of chain
1	L	433	 69% 15% 16%
2	A	556	 83% 14% .
2	B	556	 84% 13% .
2	E	556	 84% 13% .
2	G	556	 85% 13% .
2	I	556	 84% 14% .
2	K	556	 83% 14% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 41970 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 Protein TOO MANY MOUTHS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	365	Total	C	N	O	S	0	0	0
			2840	1792	513	524	11			
1	F	365	Total	C	N	O	S	0	0	0
			2840	1792	513	524	11			
1	D	365	Total	C	N	O	S	0	0	0
			2840	1792	513	524	11			
1	L	365	Total	C	N	O	S	0	0	0
			2840	1792	513	524	11			
1	H	365	Total	C	N	O	S	0	0	0
			2840	1792	513	524	11			
1	J	365	Total	C	N	O	S	0	0	0
			2840	1792	513	524	11			

- Molecule 2 is a protein called LRR receptor-like serine/threonine-protein kinase ERL1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	544	Total	C	N	O	S	0	0	0
			4155	2641	705	791	18			
2	G	544	Total	C	N	O	S	0	0	0
			4155	2641	705	791	18			
2	A	544	Total	C	N	O	S	0	0	0
			4155	2641	705	791	18			
2	I	544	Total	C	N	O	S	0	0	0
			4155	2641	705	791	18			
2	B	544	Total	C	N	O	S	0	0	0
			4155	2641	705	791	18			
2	E	544	Total	C	N	O	S	0	0	0
			4155	2641	705	791	18			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	574	HIS	-	expression tag	UNP C0LGW6

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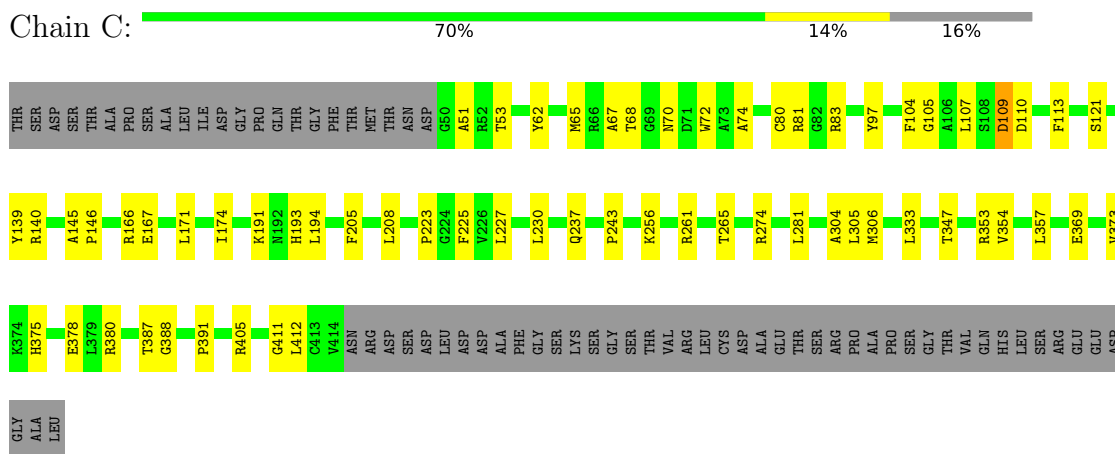
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Chain	Residue	Modelled	Actual	Comment	Reference
K	575	HIS	-	expression tag	UNP C0LGW6
K	576	HIS	-	expression tag	UNP C0LGW6
K	577	HIS	-	expression tag	UNP C0LGW6
K	578	HIS	-	expression tag	UNP C0LGW6
K	579	HIS	-	expression tag	UNP C0LGW6
G	574	HIS	-	expression tag	UNP C0LGW6
G	575	HIS	-	expression tag	UNP C0LGW6
G	576	HIS	-	expression tag	UNP C0LGW6
G	577	HIS	-	expression tag	UNP C0LGW6
G	578	HIS	-	expression tag	UNP C0LGW6
G	579	HIS	-	expression tag	UNP C0LGW6
A	574	HIS	-	expression tag	UNP C0LGW6
A	575	HIS	-	expression tag	UNP C0LGW6
A	576	HIS	-	expression tag	UNP C0LGW6
A	577	HIS	-	expression tag	UNP C0LGW6
A	578	HIS	-	expression tag	UNP C0LGW6
A	579	HIS	-	expression tag	UNP C0LGW6
I	574	HIS	-	expression tag	UNP C0LGW6
I	575	HIS	-	expression tag	UNP C0LGW6
I	576	HIS	-	expression tag	UNP C0LGW6
I	577	HIS	-	expression tag	UNP C0LGW6
I	578	HIS	-	expression tag	UNP C0LGW6
I	579	HIS	-	expression tag	UNP C0LGW6
B	574	HIS	-	expression tag	UNP C0LGW6
B	575	HIS	-	expression tag	UNP C0LGW6
B	576	HIS	-	expression tag	UNP C0LGW6
B	577	HIS	-	expression tag	UNP C0LGW6
B	578	HIS	-	expression tag	UNP C0LGW6
B	579	HIS	-	expression tag	UNP C0LGW6
E	574	HIS	-	expression tag	UNP C0LGW6
E	575	HIS	-	expression tag	UNP C0LGW6
E	576	HIS	-	expression tag	UNP C0LGW6
E	577	HIS	-	expression tag	UNP C0LGW6
E	578	HIS	-	expression tag	UNP C0LGW6
E	579	HIS	-	expression tag	UNP C0LGW6

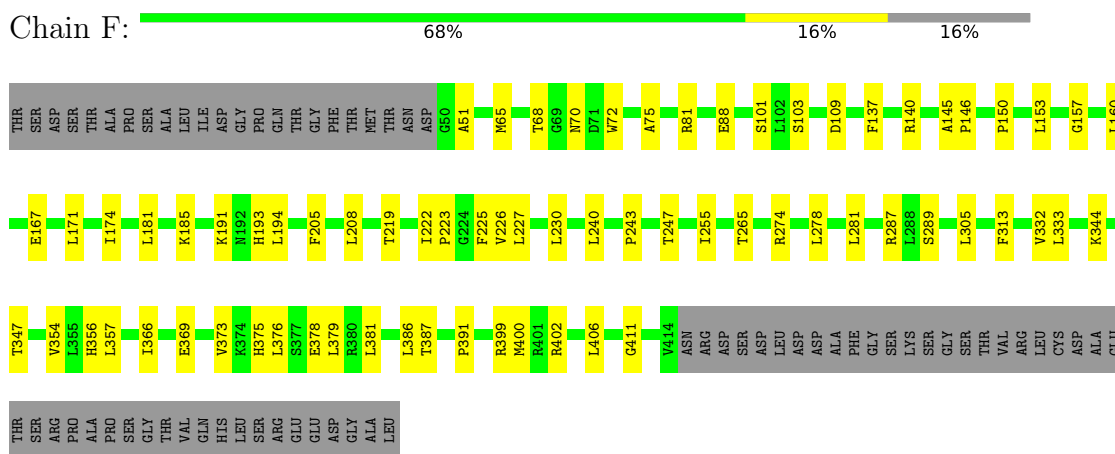
3 Residue-property plots

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.

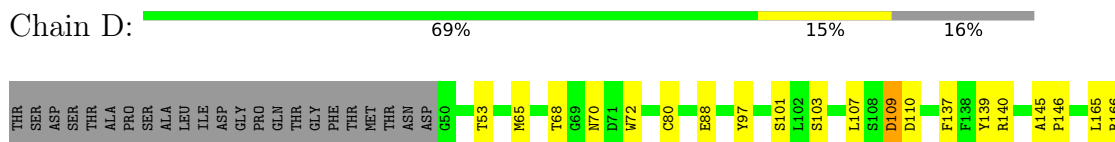
• Molecule 1: Protein TOO MANY MOUTHS

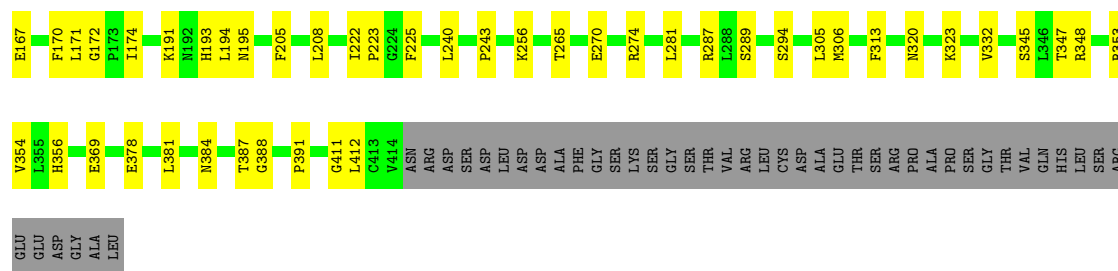


• Molecule 1: Protein TOO MANY MOUTHS



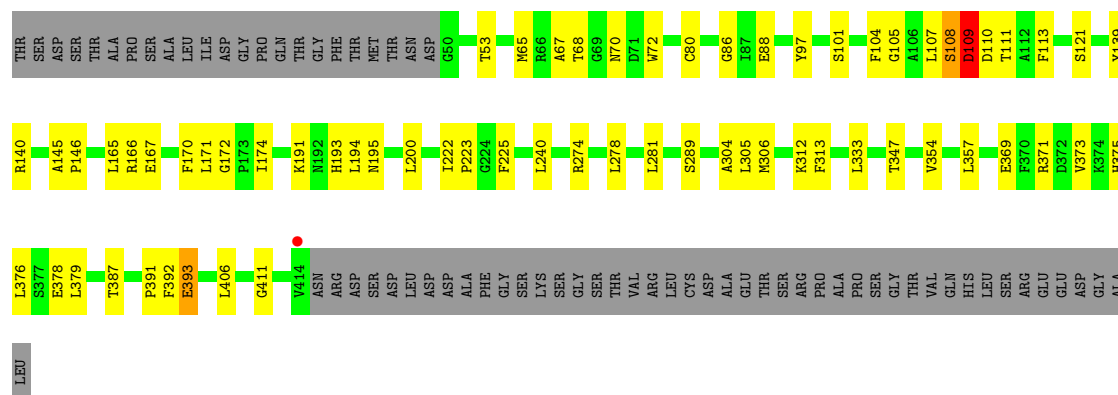
• Molecule 1: Protein TOO MANY MOUTHS





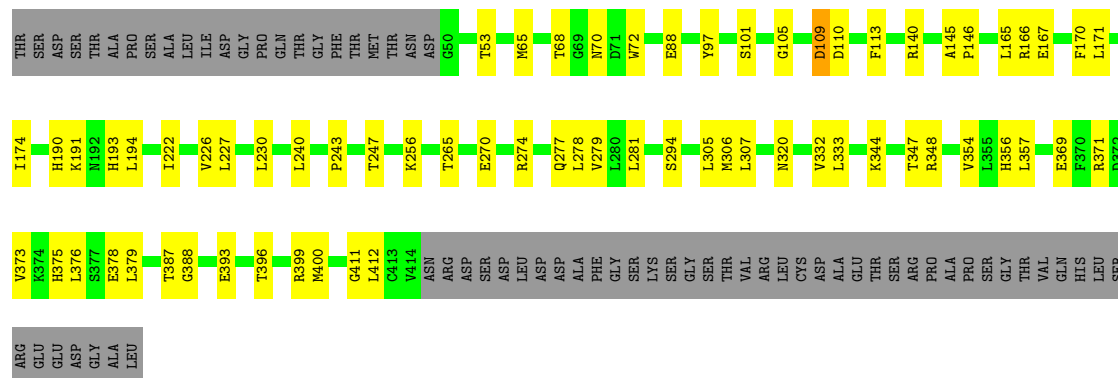
• Molecule 1: Protein TOO MANY MOUTHS

Chain L: 69% 15% 16%



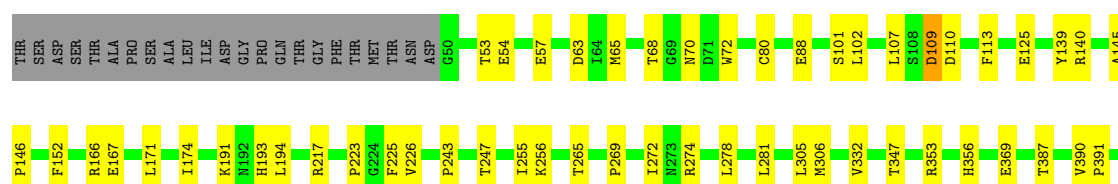
• Molecule 1: Protein TOO MANY MOUTHS

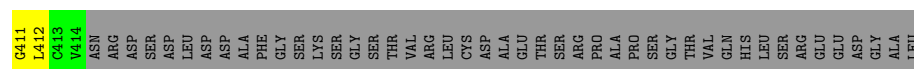
Chain H: 69% 15% 16%



• Molecule 1: Protein TOO MANY MOUTHS

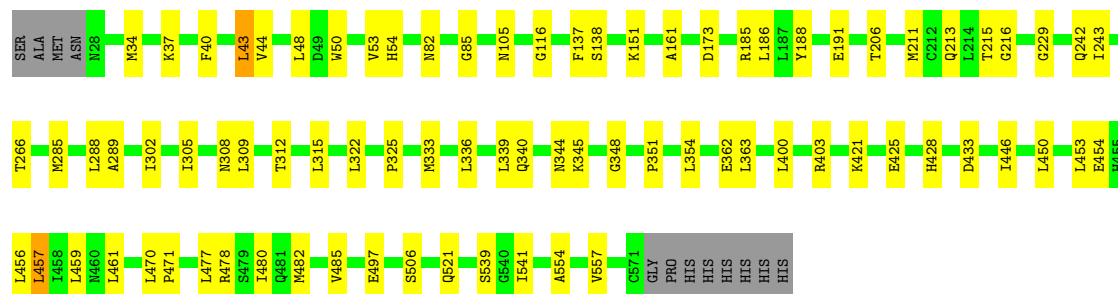
Chain J: 72% 12% 16%





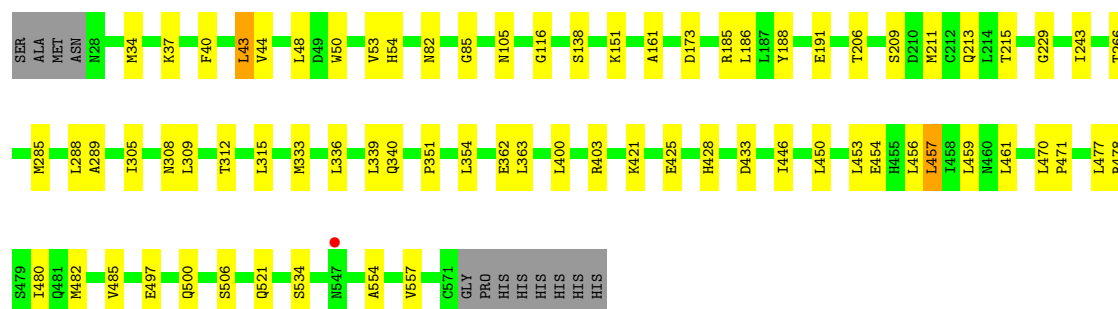
- Molecule 2: LRR receptor-like serine/threonine-protein kinase ERL1

Chain K: 83% 14%



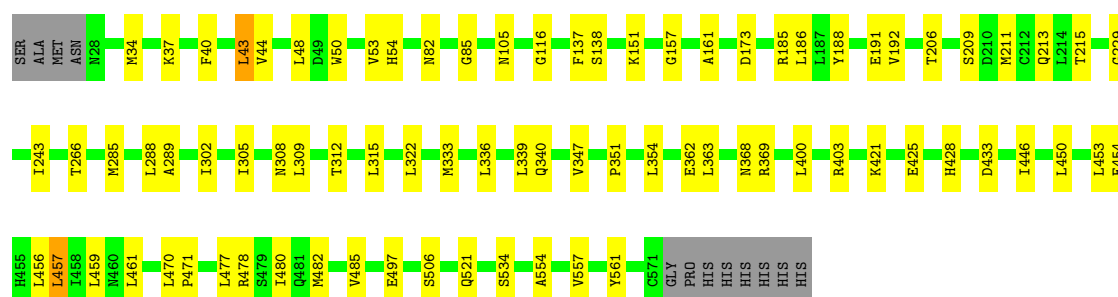
- Molecule 2: LRR receptor-like serine/threonine-protein kinase ERL1

Chain G: 85% 13%



- Molecule 2: LRR receptor-like serine/threonine-protein kinase ERL1

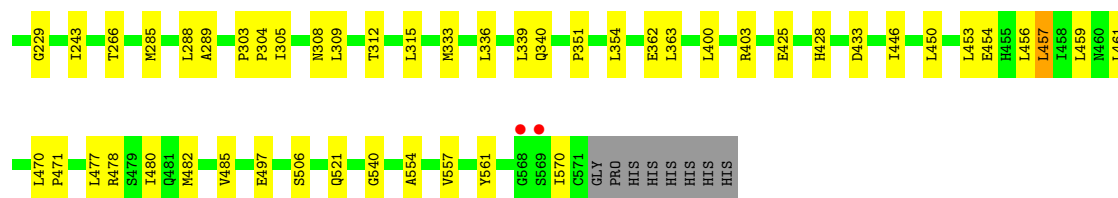
Chain A: 83% 14%



- Molecule 2: LRR receptor-like serine/threonine-protein kinase ERL1

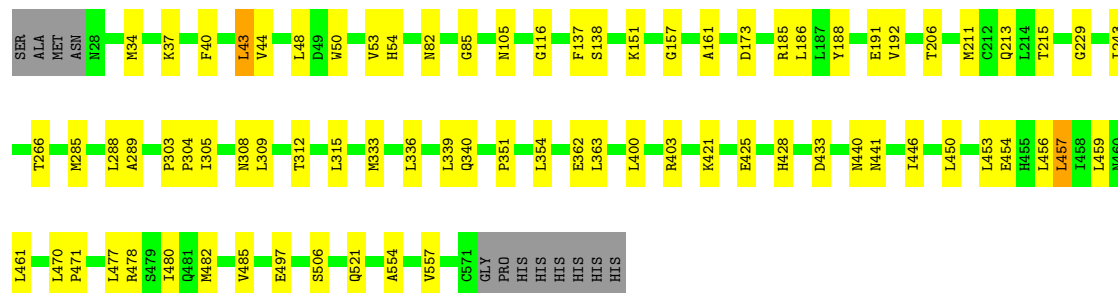
Chain I: 84% 14%





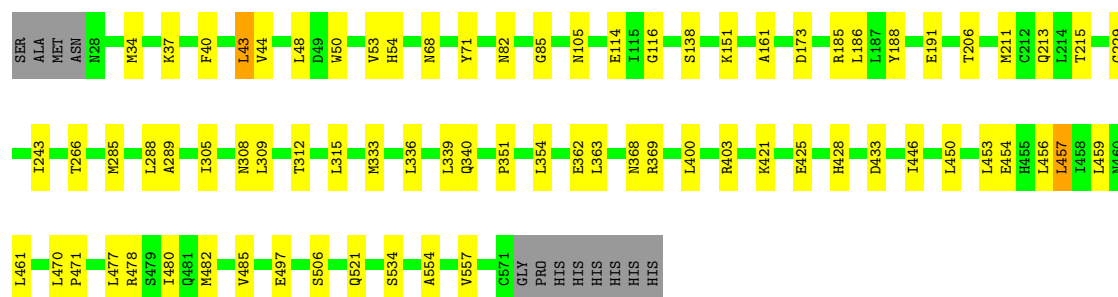
- Molecule 2: LRR receptor-like serine/threonine-protein kinase ERL1

Chain B: 84% 13%



- Molecule 2: LRR receptor-like serine/threonine-protein kinase ERL1

Chain E: 84% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	116.17Å 114.39Å 191.95Å 89.57° 90.46° 59.86°	Depositor
Resolution (Å)	43.25 – 3.06 43.25 – 3.06	Depositor EDS
% Data completeness (in resolution range)	97.6 (43.25-3.06) 97.6 (43.25-3.06)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.59 (at 3.07Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.254 , 0.276 0.282 , 0.295	Depositor DCC
R_{free} test set	7889 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	58.8	Xtriage
Anisotropy	0.499	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 26.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for k,-h+k,l 0.000 for h-k,h,l 0.011 for -h+k,-h,l 0.011 for -k,h-k,l 0.087 for -h+k,k,-l 0.016 for -k,-h,-l 0.002 for -h,-k,l 0.002 for -h,-h+k,-l 0.031 for h,h-k,-l 0.000 for h-k,-k,-l 0.001 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	41970	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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	C	0.15	0/2896	0.44	0/3933
1	D	0.15	0/2896	0.44	0/3933
1	F	0.14	0/2896	0.43	0/3933
1	H	0.26	0/2896	0.49	1/3933 (0.0%)
1	J	0.15	0/2896	0.43	0/3933
1	L	0.34	0/2896	0.53	3/3933 (0.1%)
2	A	0.15	0/4231	0.40	0/5758
2	B	0.15	0/4231	0.40	0/5758
2	E	0.15	0/4231	0.40	0/5758
2	G	0.15	0/4231	0.40	0/5758
2	I	0.15	0/4231	0.40	0/5758
2	K	0.15	0/4231	0.40	0/5758
All	All	0.18	0/42762	0.42	4/58146 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	108	SER	CA-C-N	6.78	134.49	121.54
1	L	108	SER	C-N-CA	6.78	134.49	121.54
1	H	396	THR	N-CA-C	6.02	117.93	111.36
1	L	109	ASP	N-CA-CB	5.44	119.68	110.49

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2840	0	2889	39	3
1	D	2840	0	2889	45	0
1	F	2840	0	2889	47	0
1	H	2840	0	2889	44	0
1	J	2840	0	2889	39	1
1	L	2840	0	2889	46	0
2	A	4155	0	4195	59	1
2	B	4155	0	4195	55	0
2	E	4155	0	4195	56	0
2	G	4155	0	4195	54	0
2	I	4155	0	4195	53	3
2	K	4155	0	4195	61	0
All	All	41970	0	42504	557	4

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 7.

All (557) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:242:GLN:HG3	1:L:110:ASP:OD2	1.21	1.29
2:K:242:GLN:CG	1:L:110:ASP:OD2	2.12	0.98
2:G:454:GLU:OE1	2:G:478:ARG:NH1	1.97	0.97
2:A:454:GLU:OE1	2:A:478:ARG:NH1	1.97	0.97
2:I:454:GLU:OE1	2:I:478:ARG:NH1	1.97	0.97
2:E:454:GLU:OE1	2:E:478:ARG:NH1	1.97	0.97
2:K:454:GLU:OE1	2:K:478:ARG:NH1	1.97	0.96
2:B:454:GLU:OE1	2:B:478:ARG:NH1	1.97	0.96
2:G:340:GLN:NE2	2:G:362:GLU:OE2	2.03	0.92
2:K:340:GLN:NE2	2:K:362:GLU:OE2	2.03	0.92
2:B:340:GLN:NE2	2:B:362:GLU:OE2	2.03	0.92
2:I:340:GLN:NE2	2:I:362:GLU:OE2	2.03	0.91
2:E:340:GLN:NE2	2:E:362:GLU:OE2	2.03	0.90
2:A:340:GLN:NE2	2:A:362:GLU:OE2	2.03	0.90
1:J:70:ASN:HD22	1:J:72:TRP:HE1	1.21	0.88
1:L:68:THR:HG21	1:L:104:PHE:HE1	1.39	0.87
1:L:68:THR:HG21	1:L:104:PHE:CE1	2.11	0.85
2:K:400:LEU:HD22	1:J:125:GLU:HB3	1.59	0.84
1:C:70:ASN:HD22	1:C:72:TRP:HE1	1.23	0.84
1:D:70:ASN:HD22	1:D:72:TRP:HE1	1.27	0.82
2:E:116:GLY:HA3	2:E:138:SER:HB2	1.62	0.82
2:K:116:GLY:HA3	2:K:138:SER:HB2	1.62	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:116:GLY:HA3	2:G:138:SER:HB2	1.62	0.82
2:B:116:GLY:HA3	2:B:138:SER:HB2	1.62	0.81
2:A:116:GLY:HA3	2:A:138:SER:HB2	1.62	0.81
2:I:482:MET:HG2	2:I:506:SER:HB2	1.63	0.81
2:B:482:MET:HG2	2:B:506:SER:HB2	1.63	0.81
2:I:116:GLY:HA3	2:I:138:SER:HB2	1.62	0.81
2:E:482:MET:HG2	2:E:506:SER:HB2	1.63	0.80
2:G:482:MET:HG2	2:G:506:SER:HB2	1.63	0.80
2:K:482:MET:HG2	2:K:506:SER:HB2	1.63	0.80
1:H:145:ALA:HB1	1:H:146:PRO:HD2	1.65	0.79
1:F:145:ALA:HB1	1:F:146:PRO:HD2	1.66	0.78
2:A:482:MET:HG2	2:A:506:SER:HB2	1.63	0.78
2:G:433:ASP:HB2	2:G:457:LEU:HD22	1.66	0.77
2:A:433:ASP:HB2	2:A:457:LEU:HD22	1.67	0.77
2:E:433:ASP:HB2	2:E:457:LEU:HD22	1.67	0.77
2:I:433:ASP:HB2	2:I:457:LEU:HD22	1.67	0.77
2:B:433:ASP:HB2	2:B:457:LEU:HD22	1.67	0.77
2:K:433:ASP:HB2	2:K:457:LEU:HD22	1.66	0.77
2:I:351:PRO:HG2	2:I:354:LEU:HD13	1.68	0.76
2:B:351:PRO:HG2	2:B:354:LEU:HD13	1.68	0.76
2:E:351:PRO:HG2	2:E:354:LEU:HD13	1.68	0.76
2:K:351:PRO:HG2	2:K:354:LEU:HD13	1.68	0.76
1:J:145:ALA:HB1	1:J:146:PRO:HD2	1.68	0.75
2:A:351:PRO:HG2	2:A:354:LEU:HD13	1.68	0.75
2:K:325:PRO:HB3	1:D:320:ASN:OD1	1.86	0.75
2:G:351:PRO:HG2	2:G:354:LEU:HD13	1.68	0.75
1:C:145:ALA:HB1	1:C:146:PRO:HD2	1.68	0.75
2:K:497:GLU:OE2	2:K:497:GLU:N	2.20	0.75
2:A:497:GLU:OE2	2:A:497:GLU:N	2.20	0.74
2:B:497:GLU:N	2:B:497:GLU:OE2	2.20	0.74
2:I:497:GLU:N	2:I:497:GLU:OE2	2.20	0.74
2:E:497:GLU:N	2:E:497:GLU:OE2	2.20	0.73
1:J:387:THR:HA	1:J:411:GLY:HA3	1.70	0.73
1:L:70:ASN:HD22	1:L:72:TRP:HE1	1.35	0.72
1:L:145:ALA:HB1	1:L:146:PRO:HD2	1.71	0.71
2:G:497:GLU:N	2:G:497:GLU:OE2	2.20	0.71
1:D:145:ALA:HB1	1:D:146:PRO:HD2	1.71	0.71
1:C:387:THR:HA	1:C:411:GLY:HA3	1.73	0.70
1:F:70:ASN:HD22	1:F:72:TRP:HE1	1.40	0.69
1:H:70:ASN:HD22	1:H:72:TRP:HE1	1.39	0.69
1:H:320:ASN:ND2	2:A:347:VAL:HG23	2.07	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:387:THR:HA	1:D:411:GLY:HA3	1.76	0.67
1:F:387:THR:HA	1:F:411:GLY:HA3	1.77	0.66
2:G:206:THR:HG23	2:G:229:GLY:HA3	1.79	0.65
2:K:541:ILE:HG23	1:F:75:ALA:HB1	1.79	0.65
2:G:500:GLN:NE2	2:E:521:GLN:HE21	1.94	0.65
2:A:206:THR:HG23	2:A:229:GLY:HA3	1.79	0.65
2:K:206:THR:HG23	2:K:229:GLY:HA3	1.79	0.64
2:E:206:THR:HG23	2:E:229:GLY:HA3	1.79	0.64
2:I:206:THR:HG23	2:I:229:GLY:HA3	1.79	0.63
2:B:206:THR:HG23	2:B:229:GLY:HA3	1.79	0.63
1:H:387:THR:HA	1:H:411:GLY:HA3	1.79	0.63
1:D:171:LEU:HD23	1:D:193:HIS:HB2	1.79	0.62
1:L:387:THR:HA	1:L:411:GLY:HA3	1.82	0.60
1:C:274:ARG:HG2	1:C:274:ARG:HH11	1.67	0.60
2:B:305:ILE:O	2:B:308:ASN:ND2	2.36	0.59
1:F:174:ILE:HG12	1:F:194:LEU:HD13	1.85	0.59
1:D:270:GLU:OE2	1:D:294:SER:OG	2.11	0.59
2:K:305:ILE:O	2:K:308:ASN:ND2	2.36	0.59
1:H:320:ASN:CG	2:A:347:VAL:HG23	2.28	0.59
1:L:289:SER:HA	1:L:313:PHE:HA	1.85	0.59
2:G:305:ILE:O	2:G:308:ASN:ND2	2.36	0.58
2:A:305:ILE:O	2:A:308:ASN:ND2	2.36	0.58
1:F:332:VAL:HA	1:F:356:HIS:HB2	1.85	0.58
1:C:53:THR:HB	1:C:80:CYS:HB2	1.85	0.58
2:E:305:ILE:O	2:E:308:ASN:ND2	2.36	0.58
1:L:171:LEU:HD23	1:L:193:HIS:HB2	1.85	0.58
1:H:274:ARG:HG2	1:H:274:ARG:HH11	1.69	0.58
2:I:185:ARG:HA	2:I:188:TYR:HD2	1.69	0.57
2:B:185:ARG:HA	2:B:188:TYR:HD2	1.69	0.57
2:K:185:ARG:HA	2:K:188:TYR:HD2	1.69	0.57
1:D:306:MET:HE2	2:B:137:PHE:HE1	1.70	0.57
1:J:274:ARG:HG2	1:J:274:ARG:HH11	1.68	0.57
2:A:425:GLU:HA	2:A:428:HIS:HD2	1.70	0.57
2:E:34:MET:HE2	2:E:34:MET:HA	1.87	0.57
1:H:226:VAL:HG22	1:H:247:THR:HB	1.85	0.57
2:G:185:ARG:HA	2:G:188:TYR:HD2	1.69	0.57
2:K:34:MET:HA	2:K:34:MET:HE2	1.87	0.57
1:D:347:THR:HG21	1:D:369:GLU:HB3	1.87	0.57
1:H:88:GLU:HB2	1:H:101:SER:HB3	1.86	0.57
1:F:51:ALA:O	1:F:81:ARG:NH1	2.36	0.57
2:A:53:VAL:HG13	2:A:54:HIS:ND1	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:185:ARG:HA	2:A:188:TYR:HD2	1.69	0.57
1:H:320:ASN:CG	2:A:347:VAL:CG2	2.78	0.57
1:D:88:GLU:HB2	1:D:101:SER:HB3	1.86	0.56
2:A:34:MET:HA	2:A:34:MET:HE2	1.87	0.56
2:E:185:ARG:HA	2:E:188:TYR:HD2	1.69	0.56
2:G:53:VAL:HG13	2:G:54:HIS:ND1	2.20	0.56
1:F:171:LEU:HD23	1:F:193:HIS:HB2	1.87	0.56
1:D:222:ILE:HG12	1:D:240:LEU:HD13	1.87	0.56
1:D:289:SER:HA	1:D:313:PHE:HA	1.87	0.56
2:G:425:GLU:HA	2:G:428:HIS:HD2	1.70	0.56
2:E:425:GLU:HA	2:E:428:HIS:HD2	1.70	0.56
2:I:53:VAL:HG13	2:I:54:HIS:ND1	2.20	0.56
1:C:171:LEU:HD23	1:C:193:HIS:HB2	1.86	0.56
2:A:185:ARG:HA	2:A:188:TYR:CD2	2.41	0.56
2:E:185:ARG:HA	2:E:188:TYR:CD2	2.41	0.56
1:D:274:ARG:HG2	1:D:274:ARG:HH11	1.70	0.56
2:B:425:GLU:HA	2:B:428:HIS:HD2	1.70	0.56
2:K:40:PHE:HA	2:K:85:GLY:HA3	1.88	0.56
2:G:40:PHE:HA	2:G:85:GLY:HA3	1.88	0.56
2:I:470:LEU:HD12	2:I:471:PRO:HD2	1.88	0.56
2:B:470:LEU:HD12	2:B:471:PRO:HD2	1.88	0.56
2:K:53:VAL:HG13	2:K:54:HIS:ND1	2.20	0.56
1:F:274:ARG:HG2	1:F:274:ARG:HH11	1.70	0.56
2:G:185:ARG:HA	2:G:188:TYR:CD2	2.41	0.56
2:I:425:GLU:HA	2:I:428:HIS:HD2	1.70	0.56
2:E:40:PHE:HA	2:E:85:GLY:HA3	1.88	0.56
1:D:167:GLU:HG2	1:D:191:LYS:HB3	1.88	0.56
2:G:34:MET:HE2	2:G:34:MET:HA	1.87	0.56
2:A:40:PHE:HA	2:A:85:GLY:HA3	1.88	0.56
2:B:53:VAL:HG13	2:B:54:HIS:ND1	2.20	0.56
2:B:185:ARG:HA	2:B:188:TYR:CD2	2.41	0.56
1:C:167:GLU:HG2	1:C:191:LYS:HB3	1.88	0.55
1:L:200:LEU:HD23	1:L:223:PRO:HB3	1.88	0.55
1:H:174:ILE:HG12	1:H:194:LEU:HD13	1.88	0.55
2:I:185:ARG:HA	2:I:188:TYR:CD2	2.41	0.55
2:B:40:PHE:HA	2:B:85:GLY:HA3	1.88	0.55
2:K:185:ARG:HA	2:K:188:TYR:CD2	2.41	0.55
1:H:347:THR:HG21	1:H:369:GLU:HB3	1.88	0.55
2:E:53:VAL:HG13	2:E:54:HIS:ND1	2.20	0.55
2:K:325:PRO:HG3	1:D:323:LYS:HD2	1.87	0.55
2:K:425:GLU:HA	2:K:428:HIS:HD2	1.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:470:LEU:HD12	2:E:471:PRO:HD2	1.88	0.55
1:L:274:ARG:HH11	1:L:274:ARG:HG2	1.71	0.55
2:G:470:LEU:HD12	2:G:471:PRO:HD2	1.88	0.55
2:I:40:PHE:HA	2:I:85:GLY:HA3	1.88	0.55
2:A:470:LEU:HD12	2:A:471:PRO:HD2	1.88	0.55
2:I:305:ILE:O	2:I:308:ASN:ND2	2.36	0.55
1:J:347:THR:HG21	1:J:369:GLU:HB3	1.89	0.55
2:I:34:MET:HE2	2:I:34:MET:HA	1.87	0.55
2:K:453:LEU:HD12	2:K:456:LEU:HD22	1.89	0.55
2:I:453:LEU:HD12	2:I:456:LEU:HD22	1.89	0.55
2:B:34:MET:HE2	2:B:34:MET:HA	1.87	0.55
2:K:470:LEU:HD12	2:K:471:PRO:HD2	1.88	0.54
2:G:453:LEU:HD12	2:G:456:LEU:HD22	1.89	0.54
2:B:453:LEU:HD12	2:B:456:LEU:HD22	1.89	0.54
2:K:137:PHE:HE1	1:L:306:MET:HE2	1.72	0.54
1:F:193:HIS:CE1	2:I:68:ASN:O	2.60	0.54
2:I:521:GLN:N	2:I:521:GLN:OE1	2.40	0.54
1:C:174:ILE:HG12	1:C:194:LEU:HD13	1.89	0.54
1:F:354:VAL:HG13	1:F:378:GLU:HB2	1.90	0.54
2:A:521:GLN:OE1	2:A:521:GLN:N	2.40	0.54
2:E:453:LEU:HD12	2:E:456:LEU:HD22	1.89	0.54
2:A:453:LEU:HD12	2:A:456:LEU:HD22	1.89	0.54
1:J:174:ILE:HG12	1:J:194:LEU:HD13	1.88	0.54
1:D:174:ILE:HG12	1:D:194:LEU:HD13	1.90	0.53
1:H:227:LEU:HD12	1:H:230:LEU:HD22	1.91	0.53
1:J:139:TYR:HD1	1:J:166:ARG:HB3	1.74	0.53
2:G:82:ASN:HA	2:G:105:ASN:HA	1.90	0.53
2:G:433:ASP:HA	2:G:456:LEU:HA	1.90	0.53
2:A:433:ASP:HA	2:A:456:LEU:HA	1.90	0.53
2:E:521:GLN:OE1	2:E:521:GLN:N	2.40	0.53
2:A:82:ASN:HA	2:A:105:ASN:HA	1.91	0.53
1:C:227:LEU:HD12	1:C:230:LEU:HD22	1.90	0.53
1:F:167:GLU:HG2	1:F:191:LYS:HB3	1.90	0.53
1:H:171:LEU:HD23	1:H:193:HIS:HB2	1.90	0.52
1:L:53:THR:HB	1:L:80:CYS:HB2	1.91	0.52
1:L:166:ARG:HG2	1:L:167:GLU:HG3	1.91	0.52
2:I:433:ASP:HA	2:I:456:LEU:HA	1.90	0.52
1:C:65:MET:O	1:C:68:THR:HG22	2.10	0.52
1:H:167:GLU:HG2	1:H:191:LYS:HB3	1.89	0.52
1:C:70:ASN:HB3	1:C:72:TRP:NE1	2.25	0.52
1:H:270:GLU:OE2	1:H:294:SER:OG	2.16	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:347:THR:HG21	1:C:369:GLU:HB3	1.92	0.52
1:L:88:GLU:HB2	1:L:101:SER:HB3	1.92	0.52
2:B:82:ASN:HA	2:B:105:ASN:HA	1.90	0.52
2:I:82:ASN:HA	2:I:105:ASN:HA	1.90	0.52
1:F:402:ARG:NH1	2:E:114:GLU:OE2	2.37	0.52
1:C:353:ARG:NH2	2:A:157:GLY:HA2	2.24	0.51
2:I:289:ALA:HA	2:I:312:THR:HA	1.93	0.51
1:D:139:TYR:HD1	1:D:166:ARG:HB3	1.75	0.51
2:B:433:ASP:HA	2:B:456:LEU:HA	1.90	0.51
2:E:289:ALA:HA	2:E:312:THR:HA	1.93	0.51
2:K:433:ASP:HA	2:K:456:LEU:HA	1.90	0.51
1:J:306:MET:HE2	2:I:137:PHE:HE1	1.75	0.51
2:A:289:ALA:HA	2:A:312:THR:HA	1.93	0.51
2:K:82:ASN:HA	2:K:105:ASN:HA	1.91	0.51
2:B:289:ALA:HA	2:B:312:THR:HA	1.93	0.51
2:E:433:ASP:HA	2:E:456:LEU:HA	1.90	0.51
1:C:146:PRO:HB3	1:C:171:LEU:HD12	1.92	0.51
1:J:171:LEU:HD23	1:J:193:HIS:HB2	1.92	0.51
1:H:222:ILE:HG12	1:H:240:LEU:HD13	1.93	0.51
2:E:82:ASN:HA	2:E:105:ASN:HA	1.90	0.51
2:K:289:ALA:HA	2:K:312:THR:HA	1.93	0.51
2:G:289:ALA:HA	2:G:312:THR:HA	1.93	0.51
2:K:428:HIS:NE2	1:J:63:ASP:OD1	2.44	0.51
1:L:174:ILE:HG12	1:L:194:LEU:HD13	1.92	0.51
1:J:353:ARG:NH2	2:I:157:GLY:HA2	2.26	0.51
2:K:521:GLN:OE1	2:K:521:GLN:N	2.40	0.50
1:H:399:ARG:NH2	1:H:400:MET:HE2	2.27	0.50
1:F:146:PRO:HB3	1:F:171:LEU:HD12	1.93	0.50
1:J:70:ASN:HD22	1:J:72:TRP:NE1	1.99	0.50
2:G:500:GLN:HE22	2:E:521:GLN:NE2	2.10	0.50
1:F:255:ILE:HA	1:F:278:LEU:HA	1.93	0.50
1:F:347:THR:HG21	1:F:369:GLU:HB3	1.94	0.50
1:D:107:LEU:HD21	2:B:192:VAL:HA	1.93	0.50
2:K:188:TYR:CD1	2:K:211:MET:HA	2.47	0.50
1:H:256:LYS:HB2	2:G:186:LEU:HD11	1.94	0.50
2:G:521:GLN:OE1	2:G:521:GLN:N	2.40	0.50
2:K:539:SER:HB2	1:F:72:TRP:HB3	1.94	0.50
1:F:88:GLU:HB2	1:F:101:SER:HB3	1.93	0.50
2:E:188:TYR:CD1	2:E:211:MET:HA	2.47	0.50
1:L:165:LEU:HB3	1:L:170:PHE:HE2	1.77	0.49
1:J:88:GLU:HB2	1:J:101:SER:HB3	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:34:MET:HE1	2:I:50:TRP:HB3	1.94	0.49
2:G:34:MET:HE1	2:G:50:TRP:HB3	1.95	0.49
1:J:65:MET:O	1:J:68:THR:HG22	2.12	0.49
2:A:188:TYR:CD1	2:A:211:MET:HA	2.47	0.49
1:H:281:LEU:HD23	1:H:305:LEU:HD13	1.95	0.49
2:G:188:TYR:CD1	2:G:211:MET:HA	2.47	0.49
2:I:188:TYR:CD1	2:I:211:MET:HA	2.47	0.49
2:B:34:MET:HE1	2:B:50:TRP:HB3	1.94	0.49
1:F:205:PHE:HB3	1:F:208:LEU:HB2	1.95	0.49
1:L:347:THR:HG21	1:L:369:GLU:HB3	1.95	0.49
2:A:34:MET:HE1	2:A:50:TRP:HB3	1.95	0.49
2:B:188:TYR:CD1	2:B:211:MET:HA	2.47	0.49
1:J:281:LEU:HD23	1:J:305:LEU:HD13	1.95	0.49
2:G:477:LEU:HB3	2:G:480:ILE:HB	1.95	0.48
1:C:281:LEU:HD23	1:C:305:LEU:HD13	1.95	0.48
1:D:306:MET:HE2	2:B:137:PHE:CE1	2.46	0.48
1:J:369:GLU:OE1	1:J:369:GLU:N	2.43	0.48
2:A:477:LEU:HB3	2:A:480:ILE:HB	1.95	0.48
2:B:477:LEU:HB3	2:B:480:ILE:HB	1.95	0.48
2:E:477:LEU:HB3	2:E:480:ILE:HB	1.95	0.48
1:C:369:GLU:OE1	1:C:369:GLU:N	2.41	0.48
1:F:222:ILE:HG12	1:F:240:LEU:HD13	1.95	0.48
2:I:477:LEU:HB3	2:I:480:ILE:HB	1.95	0.48
2:K:216:GLY:O	1:L:109:ASP:OD1	2.32	0.48
1:L:65:MET:O	1:L:68:THR:HG22	2.14	0.48
1:D:146:PRO:HB3	1:D:171:LEU:HD12	1.94	0.48
2:G:500:GLN:NE2	2:E:521:GLN:NE2	2.61	0.48
1:C:140:ARG:HA	1:C:167:GLU:O	2.14	0.48
2:E:34:MET:HE1	2:E:50:TRP:HB3	1.94	0.48
2:K:34:MET:HE1	2:K:50:TRP:HB3	1.95	0.48
1:L:140:ARG:HA	1:L:167:GLU:O	2.14	0.48
1:L:167:GLU:HG2	1:L:191:LYS:HB3	1.94	0.48
1:J:226:VAL:HG22	1:J:247:THR:HB	1.95	0.48
2:K:477:LEU:HB3	2:K:480:ILE:HB	1.95	0.48
1:F:376:LEU:HD23	1:F:379:LEU:HB2	1.95	0.48
1:H:369:GLU:OE1	1:H:369:GLU:N	2.44	0.48
1:H:371:ARG:HG3	1:H:393:GLU:HG3	1.96	0.48
2:B:521:GLN:OE1	2:B:521:GLN:N	2.40	0.48
1:F:157:GLY:HA2	1:F:181:LEU:HD23	1.96	0.47
2:K:461:LEU:HB2	2:K:485:VAL:HG12	1.96	0.47
1:C:306:MET:HE2	2:A:137:PHE:HE1	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:65:MET:O	1:F:68:THR:HG22	2.14	0.47
1:H:146:PRO:HB3	1:H:171:LEU:HD12	1.96	0.47
2:K:191:GLU:HA	2:K:215:THR:HG22	1.97	0.47
1:D:65:MET:O	1:D:68:THR:HG22	2.15	0.47
1:D:332:VAL:HA	1:D:356:HIS:HB2	1.96	0.47
1:H:243:PRO:HA	1:H:265:THR:O	2.14	0.47
1:J:146:PRO:HB3	1:J:171:LEU:HD12	1.96	0.47
2:E:461:LEU:HB2	2:E:485:VAL:HG12	1.96	0.47
2:E:191:GLU:HA	2:E:215:THR:HG22	1.97	0.47
1:D:53:THR:HB	1:D:80:CYS:HB2	1.97	0.47
2:A:191:GLU:HA	2:A:215:THR:HG22	1.97	0.47
2:A:461:LEU:HB2	2:A:485:VAL:HG12	1.96	0.47
2:E:34:MET:CE	2:E:50:TRP:HB3	2.45	0.47
2:K:37:LYS:HB2	2:K:50:TRP:CE3	2.50	0.47
1:L:376:LEU:HD23	1:L:379:LEU:HB2	1.97	0.47
2:K:34:MET:CE	2:K:50:TRP:HB3	2.45	0.46
2:G:161:ALA:HB1	2:G:186:LEU:HD23	1.97	0.46
1:L:371:ARG:HG3	1:L:393:GLU:CG	2.45	0.46
1:H:332:VAL:HA	1:H:356:HIS:HB2	1.97	0.46
2:G:37:LYS:HB2	2:G:50:TRP:CE3	2.50	0.46
2:A:37:LYS:HB2	2:A:50:TRP:CE3	2.51	0.46
2:I:461:LEU:HB2	2:I:485:VAL:HG12	1.96	0.46
1:J:193:HIS:CE1	2:E:68:ASN:O	2.69	0.46
2:A:34:MET:CE	2:A:50:TRP:HB3	2.45	0.46
2:E:37:LYS:HB2	2:E:50:TRP:CE3	2.50	0.46
1:D:53:THR:HA	1:D:97:TYR:HB2	1.98	0.46
2:G:34:MET:CE	2:G:50:TRP:HB3	2.45	0.46
2:I:34:MET:CE	2:I:50:TRP:HB3	2.45	0.46
2:B:37:LYS:HB2	2:B:50:TRP:CE3	2.50	0.46
2:B:461:LEU:HB2	2:B:485:VAL:HG12	1.96	0.46
2:K:348:GLY:HA3	1:D:320:ASN:CG	2.40	0.46
1:D:70:ASN:HD22	1:D:72:TRP:NE1	2.05	0.46
1:J:274:ARG:HG2	1:J:274:ARG:NH1	2.31	0.46
2:G:461:LEU:HB2	2:G:485:VAL:HG12	1.96	0.46
2:B:34:MET:CE	2:B:50:TRP:HB3	2.45	0.46
2:K:37:LYS:HD2	2:K:50:TRP:HB2	1.98	0.46
1:F:223:PRO:HB2	1:F:225:PHE:CZ	2.51	0.46
1:F:373:VAL:HG13	1:F:375:HIS:CE1	2.50	0.46
1:L:105:GLY:HA3	1:L:139:TYR:O	2.16	0.46
1:C:105:GLY:O	1:C:113:PHE:HB2	2.16	0.46
2:K:161:ALA:HB1	2:K:186:LEU:HD23	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:223:PRO:HB2	1:L:225:PHE:CZ	2.50	0.46
2:B:191:GLU:HA	2:B:215:THR:HG22	1.97	0.46
1:H:65:MET:O	1:H:68:THR:HG22	2.16	0.46
2:G:37:LYS:HD2	2:G:50:TRP:HB2	1.98	0.46
2:B:37:LYS:HE2	2:B:43:LEU:HD22	1.98	0.46
1:C:107:LEU:HD21	2:A:192:VAL:HA	1.97	0.46
1:F:140:ARG:HA	1:F:167:GLU:O	2.15	0.46
1:L:333:LEU:HB2	1:L:357:LEU:HD23	1.97	0.46
1:H:347:THR:HG22	1:H:348:ARG:HD2	1.97	0.46
2:B:161:ALA:HB1	2:B:186:LEU:HD23	1.97	0.46
2:E:37:LYS:HE2	2:E:43:LEU:HD22	1.99	0.46
2:A:37:LYS:HD2	2:A:50:TRP:HB2	1.98	0.45
2:I:37:LYS:HB2	2:I:50:TRP:CE3	2.50	0.45
2:I:161:ALA:HB1	2:I:186:LEU:HD23	1.97	0.45
2:I:400:LEU:HD12	2:I:400:LEU:HA	1.73	0.45
1:J:107:LEU:HD21	2:I:192:VAL:HA	1.97	0.45
2:I:191:GLU:HA	2:I:215:THR:HG22	1.97	0.45
1:H:354:VAL:HG13	1:H:378:GLU:HB2	1.99	0.45
2:A:161:ALA:HB1	2:A:186:LEU:HD23	1.97	0.45
2:I:37:LYS:HE2	2:I:43:LEU:HD22	1.98	0.45
2:E:37:LYS:HD2	2:E:50:TRP:HB2	1.98	0.45
2:K:37:LYS:HE2	2:K:43:LEU:HD22	1.99	0.45
1:D:205:PHE:HB3	1:D:208:LEU:HB2	1.98	0.45
2:G:151:LYS:HB2	2:G:173:ASP:OD1	2.17	0.45
1:C:53:THR:HA	1:C:97:TYR:HB2	1.98	0.45
2:K:403:ARG:HB3	2:K:425:GLU:HB3	1.99	0.45
1:F:227:LEU:HD12	1:F:230:LEU:HD22	1.97	0.45
1:L:278:LEU:HD21	1:L:281:LEU:HB2	1.98	0.45
1:H:278:LEU:HD21	1:H:281:LEU:HB2	1.97	0.45
2:E:403:ARG:HB3	2:E:425:GLU:HB3	1.99	0.45
1:F:369:GLU:OE1	1:F:369:GLU:N	2.46	0.45
1:L:65:MET:HE3	1:L:70:ASN:HB2	1.98	0.45
2:G:37:LYS:HE2	2:G:43:LEU:HD22	1.99	0.45
2:B:37:LYS:HD2	2:B:50:TRP:HB2	1.98	0.45
1:C:68:THR:HG21	1:C:104:PHE:CE1	2.51	0.45
1:F:281:LEU:HD23	1:F:305:LEU:HD13	1.99	0.45
1:H:165:LEU:HB3	1:H:170:PHE:HE2	1.82	0.45
2:A:37:LYS:HE2	2:A:43:LEU:HD22	1.99	0.45
2:I:37:LYS:HD2	2:I:50:TRP:HB2	1.98	0.45
1:F:265:THR:HG22	1:F:287:ARG:HB2	1.99	0.45
1:L:354:VAL:HG13	1:L:378:GLU:HB2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:140:ARG:HA	1:H:167:GLU:O	2.17	0.45
2:E:161:ALA:HB1	2:E:186:LEU:HD23	1.97	0.45
1:L:373:VAL:HG13	1:L:375:HIS:CE1	2.52	0.45
1:H:279:VAL:HB	2:G:186:LEU:HD22	1.99	0.45
1:H:376:LEU:HD23	1:H:379:LEU:HB2	1.99	0.45
2:G:53:VAL:HG13	2:G:54:HIS:CE1	2.52	0.45
2:G:191:GLU:HA	2:G:215:THR:HG22	1.97	0.45
2:I:151:LYS:HB2	2:I:173:ASP:OD1	2.17	0.45
2:B:400:LEU:HD12	2:B:400:LEU:HA	1.73	0.45
2:B:151:LYS:HB2	2:B:173:ASP:OD1	2.17	0.44
2:B:403:ARG:HB3	2:B:425:GLU:HB3	1.99	0.44
2:K:151:LYS:HB2	2:K:173:ASP:OD1	2.17	0.44
1:D:388:GLY:O	1:D:412:LEU:HA	2.15	0.44
1:H:105:GLY:O	1:H:113:PHE:HB2	2.18	0.44
1:H:306:MET:O	1:H:307:LEU:HD23	2.18	0.44
2:K:53:VAL:HG13	2:K:54:HIS:CE1	2.52	0.44
2:K:137:PHE:CE1	1:L:306:MET:HE2	2.51	0.44
1:D:274:ARG:HG2	1:D:274:ARG:NH1	2.32	0.44
1:D:353:ARG:NH2	2:B:157:GLY:HA2	2.31	0.44
2:G:400:LEU:HD12	2:G:400:LEU:HA	1.73	0.44
2:G:403:ARG:HB3	2:G:425:GLU:HB3	1.99	0.44
2:A:53:VAL:HG13	2:A:54:HIS:CE1	2.52	0.44
2:E:53:VAL:HG13	2:E:54:HIS:CE1	2.52	0.44
1:C:205:PHE:HB3	1:C:208:LEU:HB2	1.98	0.44
1:D:140:ARG:HA	1:D:167:GLU:O	2.17	0.44
1:J:243:PRO:HA	1:J:265:THR:O	2.17	0.44
2:A:151:LYS:HB2	2:A:173:ASP:OD1	2.17	0.44
2:A:403:ARG:HB3	2:A:425:GLU:HB3	1.99	0.44
2:E:151:LYS:HB2	2:E:173:ASP:OD1	2.17	0.44
1:C:243:PRO:HA	1:C:265:THR:O	2.17	0.44
1:D:165:LEU:HB3	1:D:170:PHE:HE2	1.82	0.44
2:I:403:ARG:HB3	2:I:425:GLU:HB3	1.99	0.44
1:J:223:PRO:HB2	1:J:225:PHE:CZ	2.53	0.44
2:K:344:ASN:HB3	2:K:345:LYS:H	1.72	0.44
1:F:150:PRO:HG2	1:F:153:LEU:HG	2.00	0.44
1:D:223:PRO:HB2	1:D:225:PHE:CZ	2.52	0.44
2:E:368:ASN:HB3	2:E:369:ARG:H	1.71	0.44
1:F:344:LYS:O	1:F:347:THR:HB	2.18	0.44
2:I:53:VAL:HG13	2:I:54:HIS:CE1	2.52	0.44
1:C:139:TYR:HD1	1:C:166:ARG:HB3	1.83	0.44
1:J:255:ILE:HA	1:J:278:LEU:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:332:VAL:HA	1:J:356:HIS:HB2	1.99	0.44
2:B:211:MET:HE2	2:B:211:MET:HB3	1.86	0.44
1:C:274:ARG:HG2	1:C:274:ARG:NH1	2.29	0.43
1:L:108:SER:HB3	1:L:109:ASP:H	1.59	0.43
1:L:369:GLU:OE1	1:L:369:GLU:N	2.45	0.43
2:G:339:LEU:HD23	2:G:363:LEU:HD13	2.00	0.43
2:K:400:LEU:HD12	2:K:400:LEU:HA	1.73	0.43
2:B:53:VAL:HG13	2:B:54:HIS:CE1	2.52	0.43
1:L:392:PHE:CE2	1:L:406:LEU:HD11	2.53	0.43
1:C:62:TYR:CD2	1:C:74:ALA:HA	2.53	0.43
1:C:380:ARG:CZ	1:C:405:ARG:NH1	2.81	0.43
1:D:223:PRO:HB2	1:D:225:PHE:CE1	2.54	0.43
1:L:146:PRO:HB3	1:L:171:LEU:HD12	2.00	0.43
1:C:109:ASP:HB3	1:C:110:ASP:H	1.70	0.43
1:F:70:ASN:HB3	1:F:72:TRP:NE1	2.33	0.43
1:H:344:LYS:O	1:H:347:THR:HB	2.19	0.43
2:B:339:LEU:HD23	2:B:363:LEU:HD13	2.00	0.43
1:D:243:PRO:HA	1:D:265:THR:O	2.18	0.43
1:L:274:ARG:HG2	1:L:274:ARG:NH1	2.32	0.43
1:J:255:ILE:HG13	2:I:186:LEU:HD13	2.00	0.43
2:K:421:LYS:HB3	2:K:421:LYS:HE2	1.81	0.43
1:D:347:THR:HG22	1:D:348:ARG:HD2	2.00	0.43
1:L:281:LEU:HD23	1:L:305:LEU:HD13	2.01	0.43
2:I:303:PRO:HA	2:I:304:PRO:HD3	1.94	0.43
1:D:265:THR:HG22	1:D:287:ARG:HB2	2.01	0.43
1:H:53:THR:HA	1:H:97:TYR:HB2	2.01	0.43
1:H:166:ARG:HG3	1:H:190:HIS:HB3	2.00	0.43
2:G:285:MET:O	2:G:288:LEU:HB2	2.19	0.43
2:G:446:ILE:HG23	2:G:450:LEU:HD22	2.01	0.43
1:C:51:ALA:O	1:C:81:ARG:NH1	2.46	0.43
2:A:339:LEU:HD23	2:A:363:LEU:HD13	2.00	0.43
2:I:339:LEU:HD23	2:I:363:LEU:HD13	2.00	0.43
2:E:285:MET:O	2:E:288:LEU:HB2	2.19	0.43
1:F:219:THR:HG22	2:I:56:SER:HB2	2.01	0.43
2:B:285:MET:O	2:B:288:LEU:HB2	2.19	0.43
1:F:160:LEU:HD23	1:F:181:LEU:HD13	2.00	0.42
1:H:277:GLN:OE1	2:G:185:ARG:NH2	2.52	0.42
2:G:336:LEU:HD21	2:G:339:LEU:HD13	2.01	0.42
2:A:400:LEU:O	2:A:403:ARG:HG2	2.19	0.42
2:A:446:ILE:HG23	2:A:450:LEU:HD22	2.01	0.42
2:E:421:LYS:H	2:E:421:LYS:HG2	1.70	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:315:LEU:HD22	2:K:333:MET:HE1	2.02	0.42
2:K:336:LEU:HD21	2:K:339:LEU:HD13	2.01	0.42
1:F:289:SER:HA	1:F:313:PHE:HA	2.02	0.42
1:F:381:LEU:HB2	1:F:406:LEU:HD23	2.02	0.42
2:A:368:ASN:HB3	2:A:369:ARG:H	1.71	0.42
2:K:285:MET:O	2:K:288:LEU:HB2	2.19	0.42
2:K:400:LEU:O	2:K:403:ARG:HG2	2.19	0.42
1:F:274:ARG:HG2	1:F:274:ARG:NH1	2.32	0.42
1:H:388:GLY:O	1:H:412:LEU:HA	2.20	0.42
2:G:400:LEU:O	2:G:403:ARG:HG2	2.19	0.42
2:I:285:MET:O	2:I:288:LEU:HB2	2.19	0.42
2:B:336:LEU:HD21	2:B:339:LEU:HD13	2.01	0.42
2:E:336:LEU:HD21	2:E:339:LEU:HD13	2.02	0.42
2:E:421:LYS:HB3	2:E:421:LYS:HE2	1.81	0.42
1:L:107:LEU:HD13	1:L:107:LEU:O	2.20	0.42
1:J:217:ARG:HD3	2:E:71:TYR:CE2	2.53	0.42
2:I:400:LEU:O	2:I:403:ARG:HG2	2.19	0.42
2:E:211:MET:HE2	2:E:211:MET:HB3	1.86	0.42
2:E:339:LEU:HD23	2:E:363:LEU:HD13	2.00	0.42
2:E:400:LEU:HA	2:E:400:LEU:HD12	1.73	0.42
1:J:269:PRO:O	1:J:272:ILE:HG12	2.20	0.42
1:J:167:GLU:HG2	1:J:191:LYS:HB3	2.01	0.42
2:A:315:LEU:HD22	2:A:333:MET:HE1	2.02	0.42
2:E:446:ILE:HG23	2:E:450:LEU:HD22	2.01	0.42
1:H:109:ASP:HB3	1:H:110:ASP:H	1.71	0.42
1:J:53:THR:HB	1:J:80:CYS:HB2	2.01	0.42
1:J:140:ARG:HA	1:J:167:GLU:O	2.20	0.42
2:G:309:LEU:HD23	2:G:312:THR:HG21	2.02	0.42
2:A:336:LEU:HD21	2:A:339:LEU:HD13	2.01	0.42
2:A:421:LYS:H	2:A:421:LYS:HG2	1.70	0.42
2:B:446:ILE:HG23	2:B:450:LEU:HD22	2.01	0.42
2:E:315:LEU:HD22	2:E:333:MET:HE1	2.02	0.42
2:K:339:LEU:HD23	2:K:363:LEU:HD13	2.00	0.42
1:F:243:PRO:HA	1:F:265:THR:O	2.18	0.42
1:J:54:GLU:HG3	1:J:57:GLU:H	1.84	0.42
1:J:256:LYS:HB2	2:I:186:LEU:HD11	2.02	0.42
2:A:285:MET:O	2:A:288:LEU:HB2	2.19	0.42
2:A:309:LEU:HD23	2:A:312:THR:HG21	2.02	0.42
2:I:336:LEU:HD21	2:I:339:LEU:HD13	2.01	0.42
2:B:421:LYS:H	2:B:421:LYS:HG2	1.70	0.42
2:B:421:LYS:HE2	2:B:421:LYS:HB3	1.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:172:GLY:C	1:L:195:ASN:HB2	2.45	0.42
2:A:400:LEU:HA	2:A:400:LEU:HD12	1.73	0.42
2:B:400:LEU:O	2:B:403:ARG:HG2	2.19	0.42
1:C:354:VAL:HG13	1:C:378:GLU:HB2	2.02	0.42
1:F:226:VAL:HG22	1:F:247:THR:HB	2.01	0.42
1:D:172:GLY:C	1:D:195:ASN:HB2	2.45	0.42
1:L:53:THR:HA	1:L:97:TYR:HB2	2.02	0.42
2:B:440:ASN:HB3	2:B:441:ASN:H	1.75	0.42
2:E:400:LEU:O	2:E:403:ARG:HG2	2.19	0.42
1:C:223:PRO:HB2	1:C:225:PHE:CZ	2.55	0.41
1:C:333:LEU:HB2	1:C:357:LEU:HD23	2.02	0.41
2:K:554:ALA:HA	2:K:557:VAL:HG23	2.02	0.41
2:I:554:ALA:HA	2:I:557:VAL:HG23	2.02	0.41
2:B:303:PRO:HA	2:B:304:PRO:HD3	1.94	0.41
2:E:554:ALA:HA	2:E:557:VAL:HG23	2.02	0.41
1:C:67:ALA:O	1:C:121:SER:HA	2.20	0.41
1:C:388:GLY:O	1:C:412:LEU:HA	2.20	0.41
2:K:211:MET:HE2	2:K:211:MET:HB3	1.86	0.41
2:K:446:ILE:HG23	2:K:450:LEU:HD22	2.01	0.41
1:F:333:LEU:HB2	1:F:357:LEU:HD23	2.02	0.41
1:J:125:GLU:HG3	1:J:152:PHE:CG	2.55	0.41
2:I:446:ILE:HG23	2:I:450:LEU:HD22	2.01	0.41
1:F:222:ILE:H	1:F:222:ILE:HG13	1.69	0.41
1:D:103:SER:HA	1:D:137:PHE:HB2	2.02	0.41
1:H:373:VAL:HG13	1:H:375:HIS:CE1	2.55	0.41
2:E:309:LEU:HD23	2:E:312:THR:HG21	2.02	0.41
2:K:456:LEU:HD21	2:K:459:LEU:HD13	2.03	0.41
1:H:274:ARG:HG2	1:H:274:ARG:NH1	2.32	0.41
2:A:554:ALA:HA	2:A:557:VAL:HG23	2.02	0.41
1:C:373:VAL:HG13	1:C:375:HIS:CE1	2.56	0.41
1:F:185:LYS:HA	1:F:208:LEU:HA	2.01	0.41
1:J:109:ASP:HB3	1:J:110:ASP:H	1.73	0.41
1:J:194:LEU:HD23	1:J:194:LEU:HA	1.90	0.41
2:I:456:LEU:HD21	2:I:459:LEU:HD13	2.03	0.41
2:B:309:LEU:HD23	2:B:312:THR:HG21	2.02	0.41
2:K:309:LEU:HD23	2:K:312:THR:HG21	2.02	0.41
1:F:366:ILE:HG12	1:F:386:LEU:HD13	2.03	0.41
1:F:399:ARG:HH22	1:F:400:MET:HE2	1.86	0.41
1:H:320:ASN:HB2	2:A:347:VAL:HG21	2.03	0.41
1:J:102:LEU:HD12	1:J:102:LEU:HA	1.91	0.41
2:E:456:LEU:HD21	2:E:459:LEU:HD13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:323:LYS:HE2	1:D:345:SER:HB3	2.03	0.41
1:C:256:LYS:HB2	2:A:186:LEU:HD11	2.02	0.41
1:F:103:SER:HA	1:F:137:PHE:HB2	2.03	0.41
2:G:421:LYS:HB3	2:G:421:LYS:HE2	1.81	0.41
2:G:534:SER:HB2	2:G:557:VAL:O	2.21	0.41
2:A:421:LYS:HB3	2:A:421:LYS:HE2	1.81	0.41
2:A:534:SER:HB2	2:A:557:VAL:O	2.21	0.41
2:B:456:LEU:HD21	2:B:459:LEU:HD13	2.03	0.41
2:B:554:ALA:HA	2:B:557:VAL:HG23	2.02	0.41
1:D:281:LEU:HD23	1:D:305:LEU:HD13	2.03	0.41
1:D:381:LEU:O	1:D:384:ASN:ND2	2.49	0.41
2:A:209:SER:C	2:A:211:MET:N	2.79	0.41
2:I:243:ILE:HG12	2:I:266:THR:HB	2.03	0.41
1:D:354:VAL:HG13	1:D:378:GLU:HB2	2.03	0.40
1:L:312:LYS:HD3	1:L:312:LYS:HA	1.91	0.40
1:H:333:LEU:HB2	1:H:357:LEU:HD23	2.03	0.40
2:G:554:ALA:HA	2:G:557:VAL:HG23	2.02	0.40
2:I:315:LEU:HD22	2:I:333:MET:HE1	2.02	0.40
2:B:243:ILE:HG12	2:B:266:THR:HB	2.03	0.40
1:C:237:GLN:HG3	1:C:261:ARG:HD2	2.04	0.40
1:C:304:ALA:HB1	1:C:306:MET:CE	2.51	0.40
2:K:243:ILE:HG12	2:K:266:THR:HB	2.03	0.40
1:L:222:ILE:HG12	1:L:240:LEU:HD13	2.03	0.40
1:L:304:ALA:HB1	1:L:306:MET:CE	2.51	0.40
2:A:243:ILE:HG12	2:A:266:THR:HB	2.03	0.40
2:I:309:LEU:HD23	2:I:312:THR:HG21	2.02	0.40
2:E:243:ILE:HG12	2:E:266:THR:HB	2.03	0.40
1:D:109:ASP:HB3	1:D:110:ASP:H	1.75	0.40
1:D:369:GLU:OE1	1:D:369:GLU:N	2.47	0.40
1:L:86:GLY:HA2	1:L:113:PHE:CD1	2.56	0.40
1:J:390:VAL:HG12	1:J:412:LEU:HD13	2.04	0.40
2:A:302:ILE:HG12	2:A:322:LEU:HD13	2.04	0.40
2:B:315:LEU:HD22	2:B:333:MET:HE1	2.02	0.40
2:K:302:ILE:HG12	2:K:322:LEU:HD13	2.04	0.40
1:D:256:LYS:HB2	2:B:186:LEU:HD11	2.04	0.40
2:G:209:SER:C	2:G:211:MET:N	2.79	0.40
2:G:315:LEU:HD22	2:G:333:MET:HE1	2.02	0.40
2:A:456:LEU:HD21	2:A:459:LEU:HD13	2.03	0.40
2:E:534:SER:HB2	2:E:557:VAL:O	2.21	0.40
1:F:194:LEU:HD23	1:F:194:LEU:HA	1.88	0.40
1:L:67:ALA:O	1:L:121:SER:HA	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:243:ILE:HG12	2:G:266:THR:HB	2.03	0.40
2:G:456:LEU:HD21	2:G:459:LEU:HD13	2.03	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:ARG:O	2:I:570:ILE:O[1_546]	1.93	0.27
1:C:113:PHE:O	2:I:561:TYR:OH[1_546]	1.96	0.24
1:C:72:TRP:CB	2:I:540:GLY:CA[1_546]	2.16	0.04
1:J:113:PHE:O	2:A:561:TYR:OH[1_554]	2.19	0.01

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	C	363/433 (84%)	334 (92%)	27 (7%)	2 (1%)	21	48
1	D	363/433 (84%)	333 (92%)	28 (8%)	2 (1%)	21	48
1	F	363/433 (84%)	333 (92%)	28 (8%)	2 (1%)	21	48
1	H	363/433 (84%)	334 (92%)	28 (8%)	1 (0%)	36	63
1	J	363/433 (84%)	332 (92%)	29 (8%)	2 (1%)	21	48
1	L	363/433 (84%)	334 (92%)	27 (7%)	2 (1%)	21	48
2	A	542/556 (98%)	485 (90%)	56 (10%)	1 (0%)	43	70
2	B	542/556 (98%)	484 (89%)	57 (10%)	1 (0%)	43	70
2	E	542/556 (98%)	485 (90%)	56 (10%)	1 (0%)	43	70
2	G	542/556 (98%)	485 (90%)	56 (10%)	1 (0%)	43	70
2	I	542/556 (98%)	485 (90%)	56 (10%)	1 (0%)	43	70
2	K	542/556 (98%)	485 (90%)	56 (10%)	1 (0%)	43	70

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	5430/5934 (92%)	4909 (90%)	504 (9%)	17 (0%)	36	63

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	109	ASP
1	F	109	ASP
1	D	109	ASP
1	L	109	ASP
1	H	109	ASP
1	J	109	ASP
1	C	391	PRO
1	F	391	PRO
1	J	391	PRO
1	D	391	PRO
2	K	44	VAL
2	G	44	VAL
2	A	44	VAL
2	I	44	VAL
2	B	44	VAL
2	E	44	VAL
1	L	391	PRO

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	C	322/378 (85%)	322 (100%)	0	100	100
1	D	322/378 (85%)	322 (100%)	0	100	100
1	F	322/378 (85%)	322 (100%)	0	100	100
1	H	322/378 (85%)	322 (100%)	0	100	100
1	J	322/378 (85%)	322 (100%)	0	100	100
1	L	322/378 (85%)	320 (99%)	2 (1%)	78	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	A	476/486 (98%)	472 (99%)	4 (1%)	73	79
2	B	476/486 (98%)	472 (99%)	4 (1%)	73	79
2	E	476/486 (98%)	472 (99%)	4 (1%)	73	79
2	G	476/486 (98%)	472 (99%)	4 (1%)	73	79
2	I	476/486 (98%)	472 (99%)	4 (1%)	73	79
2	K	476/486 (98%)	472 (99%)	4 (1%)	73	79
All	All	4788/5184 (92%)	4762 (100%)	26 (0%)	81	82

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	K	43	LEU
2	K	48	LEU
2	K	213	GLN
2	K	457	LEU
1	L	111	THR
1	L	393	GLU
2	G	43	LEU
2	G	48	LEU
2	G	213	GLN
2	G	457	LEU
2	A	43	LEU
2	A	48	LEU
2	A	213	GLN
2	A	457	LEU
2	I	43	LEU
2	I	48	LEU
2	I	213	GLN
2	I	457	LEU
2	B	43	LEU
2	B	48	LEU
2	B	213	GLN
2	B	457	LEU
2	E	43	LEU
2	E	48	LEU
2	E	213	GLN
2	E	457	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (63) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	70	ASN
1	C	132	HIS
1	C	382	ASN
2	K	82	ASN
2	K	178	HIS
2	K	237	ASN
2	K	505	ASN
2	K	510	ASN
2	K	511	ASN
1	F	132	HIS
1	F	237	GLN
1	F	327	ASN
1	F	382	ASN
1	D	70	ASN
1	D	132	HIS
1	D	327	ASN
1	L	132	HIS
1	L	203	ASN
1	H	132	HIS
1	J	70	ASN
1	J	132	HIS
2	G	82	ASN
2	G	178	HIS
2	G	237	ASN
2	G	428	HIS
2	G	500	GLN
2	G	505	ASN
2	G	510	ASN
2	G	511	ASN
2	A	82	ASN
2	A	178	HIS
2	A	237	ASN
2	A	385	ASN
2	A	428	HIS
2	A	505	ASN
2	A	510	ASN
2	A	511	ASN
2	I	55	ASN
2	I	82	ASN
2	I	178	HIS
2	I	237	ASN
2	I	385	ASN
2	I	428	HIS

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Mol	Chain	Res	Type
2	I	441	ASN
2	I	505	ASN
2	I	510	ASN
2	I	511	ASN
2	B	55	ASN
2	B	82	ASN
2	B	178	HIS
2	B	237	ASN
2	B	428	HIS
2	B	441	ASN
2	B	505	ASN
2	B	510	ASN
2	B	511	ASN
2	E	82	ASN
2	E	178	HIS
2	E	237	ASN
2	E	428	HIS
2	E	505	ASN
2	E	510	ASN
2	E	511	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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 [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	C	365/433 (84%)	-1.10	0 100 100	30, 42, 66, 85	0
1	D	365/433 (84%)	-1.00	0 100 100	31, 51, 79, 91	0
1	F	365/433 (84%)	-1.01	0 100 100	29, 51, 75, 91	0
1	H	365/433 (84%)	-0.98	0 100 100	30, 50, 84, 100	0
1	J	365/433 (84%)	-1.13	0 100 100	29, 43, 70, 90	0
1	L	365/433 (84%)	-0.77	1 (0%) 90 80	30, 61, 99, 112	0
2	A	544/556 (97%)	-0.80	0 100 100	40, 61, 84, 109	0
2	B	544/556 (97%)	-0.91	0 100 100	40, 61, 84, 109	0
2	E	544/556 (97%)	-0.87	0 100 100	40, 61, 84, 109	0
2	G	544/556 (97%)	-0.87	1 (0%) 91 83	40, 61, 84, 109	0
2	I	544/556 (97%)	-0.81	2 (0%) 88 75	40, 61, 84, 109	0
2	K	544/556 (97%)	-0.82	0 100 100	40, 61, 84, 109	0
All	All	5454/5934 (91%)	-0.91	4 (0%) 92 86	29, 58, 84, 112	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	569	SER	3.9
1	L	414	VAL	3.1
2	I	568	GLY	2.3
2	G	547	ASN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.