



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:52 AM UTC

PDB ID : 3USI / pdb_00003usi
Title : Crystal structure of LeuT bound to L-leucine in space group P2 from lipid bicelles
Authors : Wang, H.; Elferich, J.; Gouaux, E.
Deposited on : 2011-11-23
Resolution : 3.11 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

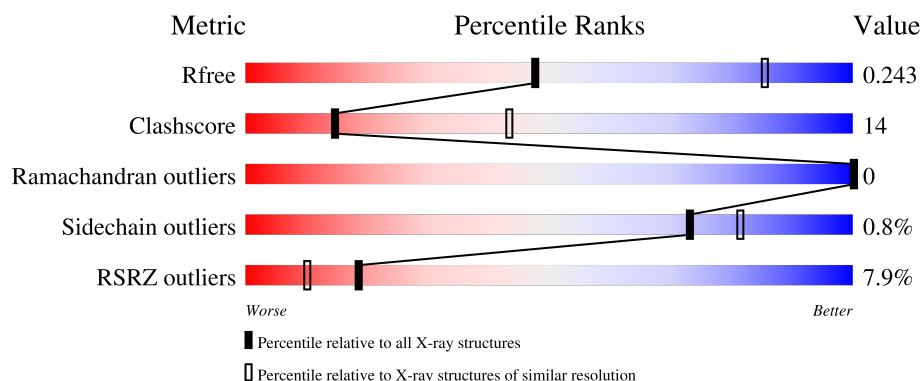
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.11 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	519	6% 67% 30%
1	B	519	9% 67% 29%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7970 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 Transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	502	Total	C	N	O	S	0	0	0
			3968	2691	622	644	11			
1	B	502	Total	C	N	O	S	0	0	0
			3968	2691	622	644	11			

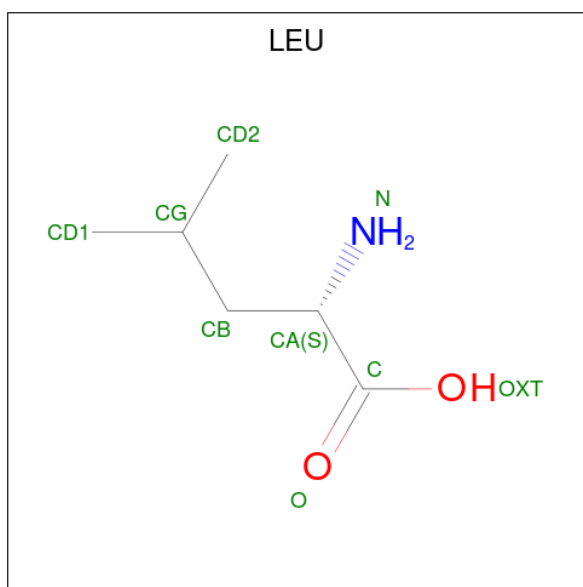
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	514	GLY	-	expression tag	UNP O67854
A	515	THR	-	expression tag	UNP O67854
A	516	LEU	-	expression tag	UNP O67854
A	517	VAL	-	expression tag	UNP O67854
A	518	PRO	-	expression tag	UNP O67854
A	519	ARG	-	expression tag	UNP O67854
B	514	GLY	-	expression tag	UNP O67854
B	515	THR	-	expression tag	UNP O67854
B	516	LEU	-	expression tag	UNP O67854
B	517	VAL	-	expression tag	UNP O67854
B	518	PRO	-	expression tag	UNP O67854
B	519	ARG	-	expression tag	UNP O67854

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Na	0	0
			2	2		
2	B	2	Total	Na	0	0
			2	2		

- Molecule 3 is LEUCINE (CCD ID: LEU) (formula: C₆H₁₃NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	6	1	2		
3	B	1	Total	C	N	O	0	0
			9	6	1	2		

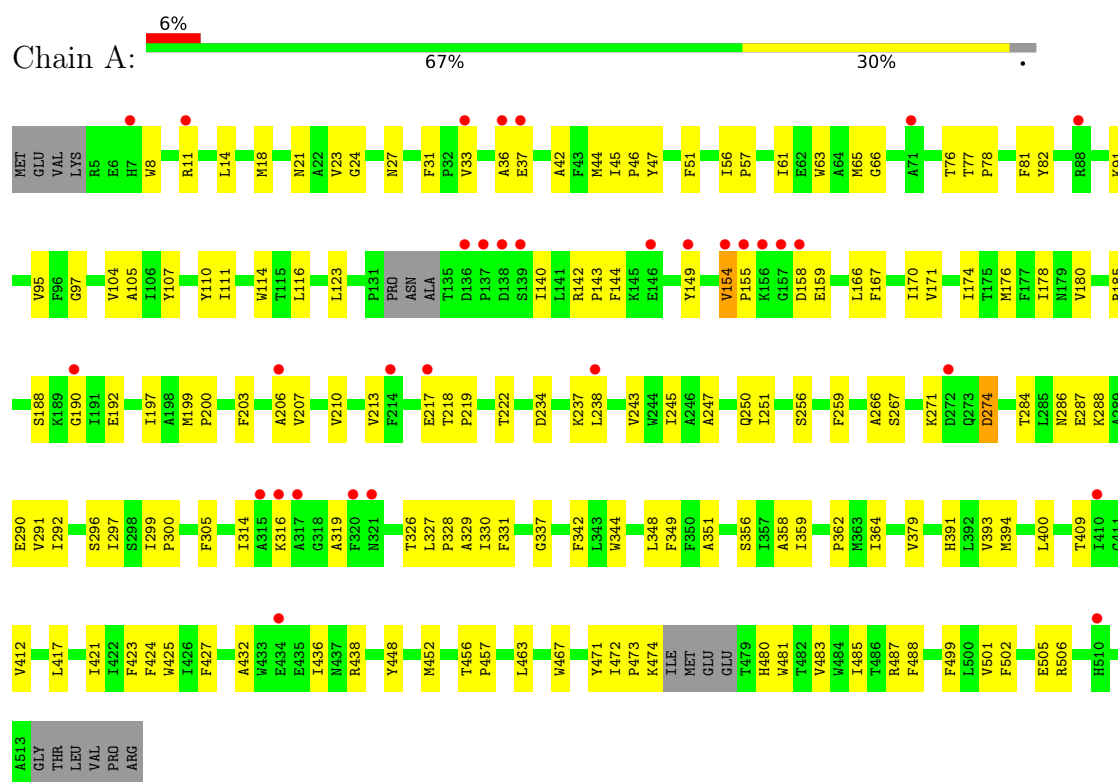
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		
4	B	6	Total	O	0	0
			6	6		

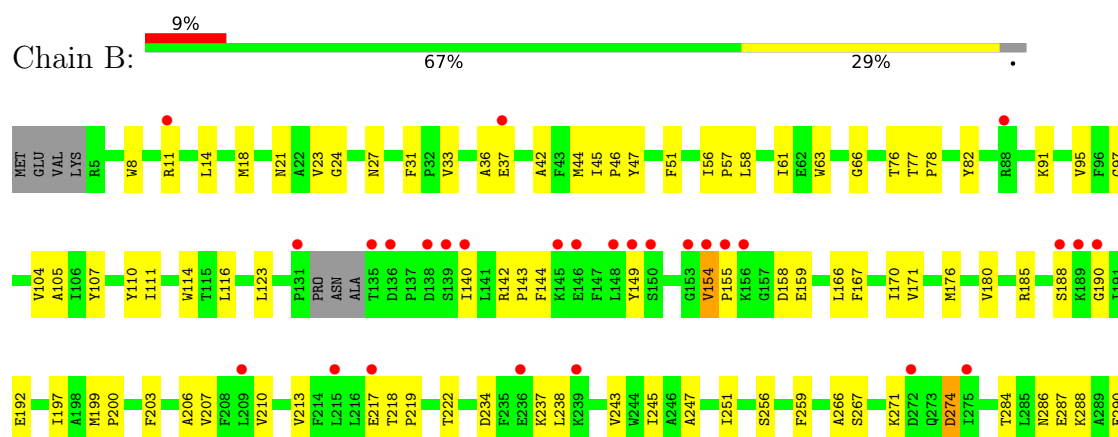
3 Residue-property plots

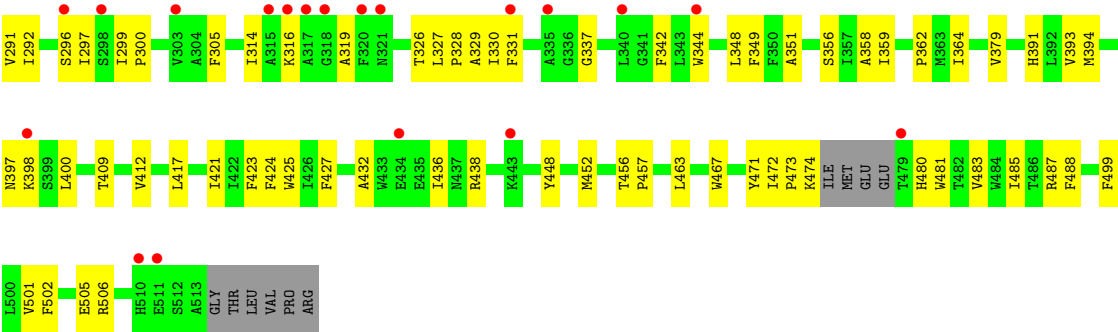
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Transporter



• Molecule 1: Transporter





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	81.18Å 84.39Å 91.03Å 90.00° 98.04° 90.00°	Depositor
Resolution (Å)	38.78 – 3.11 38.78 – 3.11	Depositor EDS
% Data completeness (in resolution range)	88.5 (38.78-3.11) 96.2 (38.78-3.11)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 3.12Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.237 , 0.250 0.239 , 0.243	Depositor DCC
R_{free} test set	1094 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	72.5	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 61.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7970	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 26.95 % 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 2.3993e-03. 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

5.1 Standard geometry

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

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	0.30	0/4093	0.72	0/5578
1	B	0.30	0/4093	0.72	0/5578
All	All	0.30	0/8186	0.72	0/11156

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3968	0	4032	112	0
1	B	3968	0	4032	111	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	9	0	10	1	0
3	B	9	0	10	2	0
4	A	6	0	0	0	0
4	B	6	0	0	0	0
All	All	7970	0	8084	223	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 14.

All (223) 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:A:24:GLY:HA2	1:A:351:ALA:HB1	1.48	0.96
1:B:33:VAL:HG21	1:B:319:ALA:HA	1.47	0.95
1:A:33:VAL:HG21	1:A:319:ALA:HA	1.47	0.92
1:B:24:GLY:HA2	1:B:351:ALA:HB1	1.53	0.89
1:A:42:ALA:HB2	1:A:234:ASP:HB3	1.57	0.85
1:B:42:ALA:HB2	1:B:234:ASP:HB3	1.57	0.85
1:B:27:ASN:HA	1:B:290:GLU:OE1	1.80	0.81
1:B:364:ILE:HD13	1:B:379:VAL:HG22	1.63	0.81
1:A:159:GLU:CD	1:A:159:GLU:H	1.90	0.80
1:A:27:ASN:HA	1:A:290:GLU:OE1	1.80	0.80
1:B:159:GLU:H	1:B:159:GLU:CD	1.90	0.78
1:A:364:ILE:HD13	1:A:379:VAL:HG22	1.63	0.78
1:B:45:ILE:HB	1:B:46:PRO:HD3	1.67	0.77
1:A:45:ILE:HB	1:A:46:PRO:HD3	1.67	0.75
1:B:481:TRP:CZ3	1:B:485:ILE:HD11	2.26	0.71
1:A:481:TRP:CZ3	1:A:485:ILE:HD11	2.25	0.71
1:A:11:ARG:HH22	1:A:274:ASP:HB3	1.58	0.68
1:B:11:ARG:HH22	1:B:274:ASP:HB3	1.58	0.67
1:B:259:PHE:CZ	1:B:412:VAL:HG11	2.32	0.65
1:A:259:PHE:CZ	1:A:412:VAL:HG11	2.32	0.65
1:A:44:MET:O	1:A:47:TYR:HB3	1.98	0.64
1:B:199:MET:HB2	1:B:200:PRO:HD3	1.81	0.63
1:A:18:MET:HG3	1:A:63:TRP:CZ2	2.34	0.63
1:B:44:MET:O	1:B:47:TYR:HB3	1.97	0.63
1:B:18:MET:HG3	1:B:63:TRP:CZ2	2.34	0.62
1:A:47:TYR:OH	1:A:290:GLU:HG2	2.00	0.61
1:B:77:THR:OG1	1:B:97:GLY:HA3	1.99	0.61
1:A:199:MET:HB2	1:A:200:PRO:HD3	1.80	0.61
1:B:47:TYR:OH	1:B:290:GLU:HG2	2.00	0.61
1:A:77:THR:OG1	1:A:97:GLY:HA3	2.00	0.61
1:B:393:VAL:HG13	1:B:400:LEU:HB2	1.83	0.61
1:A:393:VAL:HG13	1:A:400:LEU:HB2	1.83	0.60
1:A:421:ILE:O	1:A:425:TRP:HB2	2.01	0.60
1:B:421:ILE:O	1:B:425:TRP:HB2	2.01	0.59
1:A:472:ILE:N	1:A:473:PRO:HD2	2.18	0.59
1:B:472:ILE:N	1:B:473:PRO:HD2	2.18	0.58
1:B:123:LEU:HB2	1:B:342:PHE:CD1	2.38	0.58
1:B:110:TYR:CZ	1:B:394:MET:HG2	2.39	0.57
1:A:110:TYR:CZ	1:A:394:MET:HG2	2.39	0.57
1:A:123:LEU:HB2	1:A:342:PHE:CD1	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:391:HIS:HA	1:A:394:MET:HE3	1.86	0.57
1:A:331:PHE:O	1:A:337:GLY:HA3	2.05	0.57
1:B:331:PHE:O	1:B:337:GLY:HA3	2.05	0.57
1:B:296:SER:C	1:B:300:PRO:HG2	2.30	0.56
1:B:391:HIS:HA	1:B:394:MET:HE3	1.87	0.56
1:A:296:SER:C	1:A:300:PRO:HG2	2.31	0.56
1:B:432:ALA:O	1:B:436:ILE:HG13	2.06	0.55
1:A:256:SER:HG	3:A:601:LEU:N	2.05	0.55
1:A:432:ALA:O	1:A:436:ILE:HG13	2.06	0.55
1:A:327:LEU:HD23	1:A:344:TRP:CE2	2.42	0.54
1:A:111:ILE:O	1:A:114:TRP:HB2	2.08	0.54
1:A:188:SER:O	1:A:192:GLU:HB3	2.07	0.54
1:B:327:LEU:HD23	1:B:344:TRP:CE2	2.43	0.54
1:B:111:ILE:O	1:B:114:TRP:HB2	2.08	0.54
1:A:18:MET:HG3	1:A:63:TRP:CH2	2.43	0.54
1:A:154:VAL:HG13	1:A:155:PRO:HD2	1.90	0.54
1:A:158:ASP:OD1	1:A:158:ASP:C	2.50	0.54
1:A:299:ILE:HB	1:A:300:PRO:HD3	1.90	0.54
1:B:188:SER:O	1:B:192:GLU:HB3	2.07	0.54
1:B:154:VAL:HG13	1:B:155:PRO:HD2	1.90	0.53
1:B:158:ASP:C	1:B:158:ASP:OD1	2.51	0.53
1:B:66:GLY:O	1:B:267:SER:HA	2.09	0.53
1:B:18:MET:HG3	1:B:63:TRP:CH2	2.43	0.53
1:A:480:HIS:O	1:A:483:VAL:HG22	2.09	0.53
1:B:480:HIS:O	1:B:483:VAL:HG22	2.09	0.53
1:B:33:VAL:HG21	1:B:319:ALA:CA	2.31	0.53
1:A:66:GLY:O	1:A:267:SER:HA	2.09	0.52
1:A:105:ALA:HA	1:A:359:ILE:HD12	1.91	0.52
1:B:356:SER:HA	1:B:359:ILE:HG12	1.91	0.52
1:B:299:ILE:HB	1:B:300:PRO:HD3	1.90	0.52
1:B:358:ALA:O	1:B:362:PRO:HD3	2.10	0.52
1:A:31:PHE:HB2	1:A:290:GLU:OE2	2.10	0.52
1:B:217:GLU:HG3	1:B:222:THR:HG22	1.92	0.52
1:B:423:PHE:O	1:B:427:PHE:HB3	2.10	0.52
1:B:448:TYR:O	1:B:452:MET:HG3	2.10	0.52
1:A:358:ALA:O	1:A:362:PRO:HD3	2.10	0.52
1:A:423:PHE:O	1:A:427:PHE:HB3	2.10	0.51
1:B:105:ALA:HA	1:B:359:ILE:HD12	1.91	0.51
1:B:23:VAL:HG22	1:B:286:ASN:HA	1.92	0.51
1:B:256:SER:HG	3:B:601:LEU:N	2.07	0.51
1:B:284:THR:HG22	1:B:288:LYS:HE3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:GLU:HG3	1:A:222:THR:HG22	1.91	0.51
1:A:448:TYR:O	1:A:452:MET:HG3	2.10	0.51
1:B:31:PHE:HB2	1:B:290:GLU:OE2	2.10	0.51
1:A:356:SER:HA	1:A:359:ILE:HG12	1.91	0.51
1:A:142:ARG:HB3	1:A:143:PRO:HD3	1.93	0.51
1:B:159:GLU:CD	1:B:159:GLU:N	2.67	0.50
1:A:23:VAL:HG22	1:A:286:ASN:HA	1.92	0.50
1:A:284:THR:HG22	1:A:288:LYS:HE3	1.93	0.50
1:B:11:ARG:HH22	1:B:274:ASP:CB	2.24	0.50
1:A:51:PHE:CZ	1:A:56:ILE:HD11	2.47	0.49
1:B:51:PHE:CZ	1:B:56:ILE:HD11	2.47	0.49
1:A:36:ALA:HB2	1:A:299:ILE:HD11	1.94	0.49
1:A:505:GLU:HG3	1:A:506:ARG:NH1	2.28	0.49
1:B:142:ARG:HB3	1:B:143:PRO:HD3	1.94	0.49
1:B:95:VAL:HG22	1:B:501:VAL:HG21	1.95	0.49
1:B:82:TYR:CD1	1:B:91:LYS:HE3	2.48	0.49
1:A:82:TYR:CD1	1:A:91:LYS:HE3	2.48	0.48
1:B:66:GLY:HA3	1:B:266:ALA:HB3	1.95	0.48
1:B:149:TYR:CE1	1:B:154:VAL:HG21	2.48	0.48
1:A:149:TYR:CE1	1:A:154:VAL:HG21	2.48	0.48
1:A:11:ARG:HH22	1:A:274:ASP:CB	2.24	0.48
1:B:505:GLU:HG3	1:B:506:ARG:NH1	2.28	0.48
1:A:287:GLU:O	1:A:291:VAL:HG22	2.13	0.48
1:A:400:LEU:HD23	1:A:400:LEU:C	2.39	0.48
1:B:36:ALA:HB2	1:B:299:ILE:HD11	1.94	0.48
1:B:344:TRP:CE2	1:B:348:LEU:HD11	2.48	0.48
1:B:456:THR:N	1:B:457:PRO:HD2	2.28	0.48
1:A:456:THR:N	1:A:457:PRO:HD2	2.28	0.48
1:B:287:GLU:O	1:B:291:VAL:HG22	2.14	0.48
1:A:11:ARG:NH2	1:A:274:ASP:HB3	2.28	0.48
1:A:76:THR:HB	1:A:78:PRO:HD2	1.96	0.48
1:A:344:TRP:CE2	1:A:348:LEU:HD11	2.49	0.48
1:A:18:MET:O	1:A:21:ASN:HB3	2.14	0.47
1:B:18:MET:O	1:B:21:ASN:HB3	2.15	0.47
1:B:218:THR:HB	1:B:219:PRO:HD2	1.96	0.47
1:A:95:VAL:HG22	1:A:501:VAL:HG21	1.95	0.47
1:B:247:ALA:O	1:B:251:ILE:HG12	2.14	0.47
1:A:218:THR:HB	1:A:219:PRO:HD2	1.96	0.47
1:B:76:THR:HB	1:B:78:PRO:HD2	1.96	0.47
1:B:467:TRP:HD1	1:B:471:TYR:HB2	1.79	0.47
1:A:66:GLY:HA3	1:A:266:ALA:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:344:TRP:CZ2	1:B:348:LEU:HD11	2.50	0.47
1:B:400:LEU:C	1:B:400:LEU:HD23	2.39	0.47
1:A:467:TRP:HD1	1:A:471:TYR:HB2	1.79	0.47
1:B:144:PHE:CE2	1:B:329:ALA:HA	2.50	0.47
1:A:327:LEU:N	1:A:328:PRO:HD2	2.30	0.46
1:A:247:ALA:O	1:A:251:ILE:HG12	2.14	0.46
1:A:33:VAL:HG21	1:A:319:ALA:CA	2.31	0.46
1:A:344:TRP:CZ2	1:A:348:LEU:HD11	2.51	0.46
1:A:197:ILE:O	1:A:200:PRO:HD2	2.16	0.46
1:A:144:PHE:CE2	1:A:329:ALA:HA	2.51	0.46
1:B:327:LEU:N	1:B:328:PRO:HD2	2.30	0.46
1:B:57:PRO:O	1:B:61:ILE:HG13	2.16	0.46
1:B:176:MET:O	1:B:180:VAL:HG23	2.16	0.46
1:A:159:GLU:CD	1:A:159:GLU:N	2.67	0.45
1:B:45:ILE:HB	1:B:46:PRO:CD	2.43	0.45
1:A:176:MET:O	1:A:180:VAL:HG23	2.16	0.45
1:A:506:ARG:HA	1:A:506:ARG:HD3	1.76	0.45
1:A:499:PHE:O	1:A:502:PHE:HB3	2.16	0.45
1:A:291:VAL:HG23	1:A:292:ILE:HG13	1.99	0.45
1:B:56:ILE:N	1:B:57:PRO:HD2	2.32	0.45
1:A:104:VAL:HA	1:A:107:TYR:CE2	2.51	0.45
1:B:197:ILE:O	1:B:200:PRO:HD2	2.16	0.45
1:A:56:ILE:N	1:A:57:PRO:HD2	2.32	0.45
1:A:57:PRO:O	1:A:61:ILE:HG13	2.15	0.45
1:A:409:THR:HG22	1:A:467:TRP:CH2	2.51	0.45
1:B:409:THR:HG22	1:B:467:TRP:CH2	2.52	0.45
1:B:417:LEU:O	1:B:421:ILE:HG13	2.17	0.45
1:B:238:LEU:HD23	1:B:238:LEU:N	2.32	0.45
1:B:37:GLU:HG2	1:B:316:LYS:HA	1.99	0.44
1:B:499:PHE:O	1:B:502:PHE:HB3	2.16	0.44
1:A:417:LEU:O	1:A:421:ILE:HG13	2.16	0.44
1:A:238:LEU:N	1:A:238:LEU:HD23	2.32	0.44
1:A:82:TYR:CE1	1:A:91:LYS:HE3	2.53	0.44
1:A:37:GLU:HG2	1:A:316:LYS:HA	1.99	0.44
1:A:305:PHE:CE1	1:A:330:ILE:HG23	2.53	0.44
1:B:104:VAL:HA	1:B:107:TYR:CE2	2.52	0.44
1:A:154:VAL:HG13	1:A:155:PRO:CD	2.48	0.44
1:B:291:VAL:HG23	1:B:292:ILE:HG13	1.99	0.44
1:B:506:ARG:HA	1:B:506:ARG:HD3	1.76	0.44
1:B:42:ALA:HB1	1:B:237:LYS:HB2	2.00	0.43
1:B:166:LEU:O	1:B:170:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:LYS:O	1:B:292:ILE:HB	2.18	0.43
1:A:166:LEU:O	1:A:170:ILE:HG13	2.18	0.43
1:A:245:ILE:HG12	1:A:463:LEU:HD11	2.01	0.43
1:B:167:PHE:O	1:B:171:VAL:HG23	2.19	0.43
1:B:421:ILE:HG23	1:B:425:TRP:CE3	2.54	0.43
1:A:8:TRP:CD1	1:A:14:LEU:HD13	2.53	0.43
1:B:8:TRP:CD1	1:B:14:LEU:HD13	2.54	0.43
1:B:82:TYR:CE1	1:B:91:LYS:HE3	2.53	0.43
1:A:481:TRP:HZ3	1:A:485:ILE:HD11	1.80	0.43
1:B:259:PHE:CZ	3:B:601:LEU:HD12	2.54	0.43
1:A:288:LYS:O	1:A:292:ILE:HB	2.18	0.43
1:B:305:PHE:CE1	1:B:330:ILE:HG23	2.53	0.43
1:A:421:ILE:HG23	1:A:425:TRP:CE3	2.54	0.42
1:A:487:ARG:O	1:A:488:PHE:C	2.63	0.42
1:B:57:PRO:HA	1:B:452:MET:HE2	2.01	0.42
1:B:237:LYS:HE3	1:B:243:VAL:HG21	2.02	0.42
1:A:167:PHE:O	1:A:171:VAL:HG23	2.19	0.42
1:B:154:VAL:HG13	1:B:155:PRO:CD	2.48	0.42
1:A:502:PHE:O	1:A:505:GLU:HG2	2.20	0.42
1:A:271:LYS:HG2	1:A:438:ARG:HH12	1.85	0.42
1:A:213:VAL:HG21	1:A:297:ILE:HG23	2.02	0.42
1:A:237:LYS:HE3	1:A:243:VAL:HG21	2.02	0.42
1:A:359:ILE:C	1:A:362:PRO:HD2	2.45	0.42
1:B:480:HIS:CE1	1:B:481:TRP:CD1	3.08	0.42
1:A:203:PHE:O	1:A:207:VAL:HG23	2.20	0.42
1:B:58:LEU:HD23	1:B:58:LEU:HA	1.94	0.42
1:A:42:ALA:HB1	1:A:237:LYS:HB2	2.00	0.41
1:B:245:ILE:HG12	1:B:463:LEU:HD11	2.00	0.41
1:A:45:ILE:HB	1:A:46:PRO:CD	2.43	0.41
1:B:218:THR:HB	1:B:219:PRO:CD	2.50	0.41
1:A:116:LEU:HB2	1:A:349:PHE:CD1	2.56	0.41
1:B:271:LYS:HG2	1:B:438:ARG:HH12	1.85	0.41
1:A:140:ILE:HG22	1:A:329:ALA:HB1	2.01	0.41
1:B:397:ASN:O	1:B:398:LYS:HB2	2.21	0.41
1:A:471:TYR:O	1:A:474:LYS:N	2.54	0.41
1:B:356:SER:HA	1:B:359:ILE:CG1	2.51	0.41
1:B:487:ARG:O	1:B:488:PHE:C	2.62	0.41
1:A:356:SER:HA	1:A:359:ILE:CG1	2.50	0.41
1:B:140:ILE:HG22	1:B:329:ALA:HB1	2.01	0.41
1:B:203:PHE:O	1:B:207:VAL:HG23	2.20	0.41
1:A:65:MET:HE1	1:A:81:PHE:HE1	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:THR:HB	1:A:219:PRO:CD	2.50	0.41
1:B:116:LEU:HB2	1:B:349:PHE:CD1	2.56	0.41
1:B:213:VAL:HG21	1:B:297:ILE:HG23	2.02	0.41
1:B:502:PHE:O	1:B:505:GLU:HG2	2.20	0.41
1:A:11:ARG:NH2	1:A:274:ASP:OD2	2.54	0.41
1:A:185:ARG:HB2	1:A:190:GLY:CA	2.51	0.41
1:A:206:ALA:O	1:A:210:VAL:HG23	2.21	0.41
1:B:11:ARG:NH2	1:B:274:ASP:OD2	2.54	0.41
1:B:206:ALA:O	1:B:210:VAL:HG23	2.21	0.41
1:B:359:ILE:C	1:B:362:PRO:HD2	2.45	0.41
1:A:174:ILE:O	1:A:178:ILE:HG13	2.22	0.41
1:A:480:HIS:CE1	1:A:481:TRP:CD1	3.08	0.41
1:A:250:GLN:NE2	1:A:290:GLU:OE2	2.54	0.40
1:A:57:PRO:HA	1:A:452:MET:HE2	2.02	0.40
1:A:314:ILE:HG21	1:A:326:THR:OG1	2.22	0.40
1:B:471:TYR:O	1:B:474:LYS:N	2.54	0.40
1:B:185:ARG:HB2	1:B:190:GLY:CA	2.51	0.40
1:B:314:ILE:HG21	1:B:326:THR:OG1	2.22	0.40
1:B:11:ARG:NH2	1:B:274:ASP:HB3	2.29	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	496/519 (96%)	480 (97%)	16 (3%)	0	100	100
1	B	496/519 (96%)	480 (97%)	16 (3%)	0	100	100
All	All	992/1038 (96%)	960 (97%)	32 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	399/420 (95%)	396 (99%)	3 (1%)	73	81
1	B	399/420 (95%)	396 (99%)	3 (1%)	73	81
All	All	798/840 (95%)	792 (99%)	6 (1%)	73	81

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

Mol	Chain	Res	Type
1	A	154	VAL
1	A	274	ASP
1	A	424	PHE
1	B	154	VAL
1	B	274	ASP
1	B	424	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	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

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	LEU	A	601	2	6,8,8	0.94	1 (16%)	5,10,10	1.18	1 (20%)
3	LEU	B	601	2	6,8,8	0.94	1 (16%)	5,10,10	1.20	1 (20%)

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
3	LEU	A	601	2	-	0/8/8/8	-
3	LEU	B	601	2	-	0/8/8/8	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	LEU	OXT-C	-2.16	1.23	1.30
3	B	601	LEU	OXT-C	-2.16	1.23	1.30

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	601	LEU	OXT-C-O	-2.68	118.01	124.08
3	A	601	LEU	OXT-C-O	-2.64	118.09	124.08

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	LEU	1	0
3	B	601	LEU	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	502/519 (96%)	0.51	32 (6%) 25 14	28, 60, 108, 174	0
1	B	502/519 (96%)	0.67	47 (9%) 14 8	31, 68, 127, 214	0
All	All	1004/1038 (96%)	0.59	79 (7%) 18 10	28, 64, 121, 214	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	317	ALA	6.5
1	B	317	ALA	6.0
1	B	149	TYR	5.9
1	B	155	PRO	5.1
1	B	321	ASN	4.9
1	B	154	VAL	4.3
1	A	217	GLU	4.2
1	A	154	VAL	4.1
1	B	140	ILE	4.0
1	A	146	GLU	4.0
1	A	37	GLU	3.8
1	A	33	VAL	3.7
1	B	136	ASP	3.6
1	A	320	PHE	3.6
1	A	157	GLY	3.5
1	A	158	ASP	3.4
1	B	188	SER	3.3
1	A	149	TYR	3.2
1	A	272	ASP	3.2
1	B	344	TRP	3.2
1	B	318	GLY	3.2
1	B	217	GLU	3.1
1	A	136	ASP	3.0
1	B	209	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	510	HIS	2.9
1	B	316	LYS	2.9
1	A	316	LYS	2.8
1	B	320	PHE	2.8
1	A	190	GLY	2.7
1	B	272	ASP	2.7
1	B	335	ALA	2.7
1	A	155	PRO	2.6
1	B	215	LEU	2.6
1	B	146	GLU	2.6
1	B	156	LYS	2.5
1	B	37	GLU	2.5
1	A	36	ALA	2.5
1	A	156	LYS	2.4
1	B	139	SER	2.4
1	B	434	GLU	2.4
1	B	239	LYS	2.4
1	A	7	HIS	2.3
1	A	139	SER	2.3
1	A	11	ARG	2.3
1	B	138	ASP	2.3
1	B	190	GLY	2.3
1	B	315	ALA	2.3
1	B	398	LYS	2.3
1	A	315	ALA	2.2
1	B	145	LYS	2.2
1	A	88	ARG	2.2
1	A	71	ALA	2.2
1	A	238	LEU	2.2
1	B	189	LYS	2.2
1	B	88	ARG	2.2
1	B	275	ILE	2.2
1	B	479	THR	2.2
1	B	11	ARG	2.1
1	B	131	PRO	2.1
1	B	153	GLY	2.1
1	B	511	GLU	2.1
1	A	138	ASP	2.1
1	B	340	LEU	2.1
1	A	137	PRO	2.1
1	B	296	SER	2.1
1	B	135	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	214	PHE	2.1
1	B	298	SER	2.1
1	A	510	HIS	2.1
1	B	331	PHE	2.1
1	B	303	VAL	2.0
1	B	150	SER	2.0
1	A	206	ALA	2.0
1	A	434	GLU	2.0
1	B	443	LYS	2.0
1	A	321	ASN	2.0
1	B	236	GLU	2.0
1	B	148	LEU	2.0
1	A	410	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NA	B	752	1/1	0.87	0.08	49,49,49,49	0
3	LEU	A	601	9/9	0.89	0.17	53,54,56,56	0
2	NA	A	751	1/1	0.90	0.11	49,49,49,49	0
3	LEU	B	601	9/9	0.91	0.13	49,52,55,55	0
2	NA	A	752	1/1	0.94	0.08	72,72,72,72	0
2	NA	B	751	1/1	0.97	0.11	62,62,62,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.