



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:23 PM GMT

PDB ID : 3TT3
Title : Crystal Structure of LeuT in the inward-open conformation in complex with Fab
Authors : Krishnamurthy, H.; Gouaux, E.
Deposited on : 2011-09-13
Resolution : 3.22 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

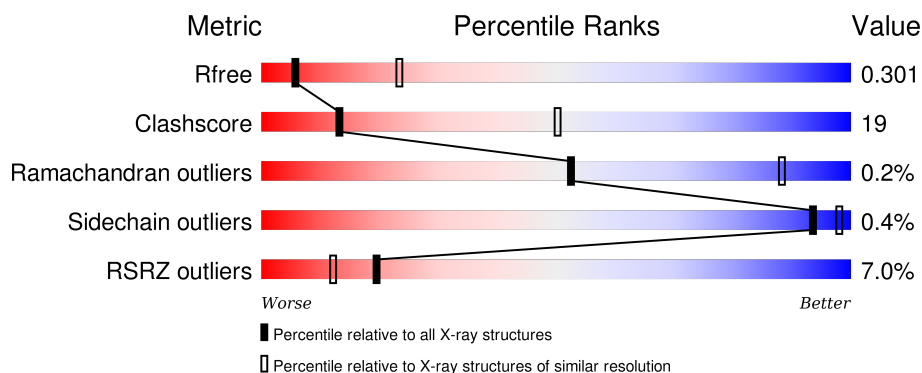
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.22 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}	91344	1095 (3.26-3.18)
Clashscore	102246	1046 (3.24-3.20)
Ramachandran outliers	100387	1026 (3.24-3.20)
Sidechain outliers	100360	1025 (3.24-3.20)
RSRZ outliers	91569	1100 (3.26-3.18)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>4%</div> <div>59%</div> <div>38%</div> <div>.</div> </div>
2	L	218	<div> <div>11%</div> <div>62%</div> <div>36%</div> <div>.</div> </div>
3	H	221	<div> <div>8%</div> <div>73%</div> <div>26%</div> <div>.</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SOG	A	520	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7200 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 Leucine transporter LeuT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	0	0	0
			3855	2607	596	641	11			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	268	ALA	TYR	ENGINEERED MUTATION	UNP O67854
A	288	ALA	LYS	ENGINEERED MUTATION	UNP O67854
A	354	VAL	THR	ENGINEERED MUTATION	UNP O67854
A	355	ALA	SER	ENGINEERED MUTATION	UNP O67854
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

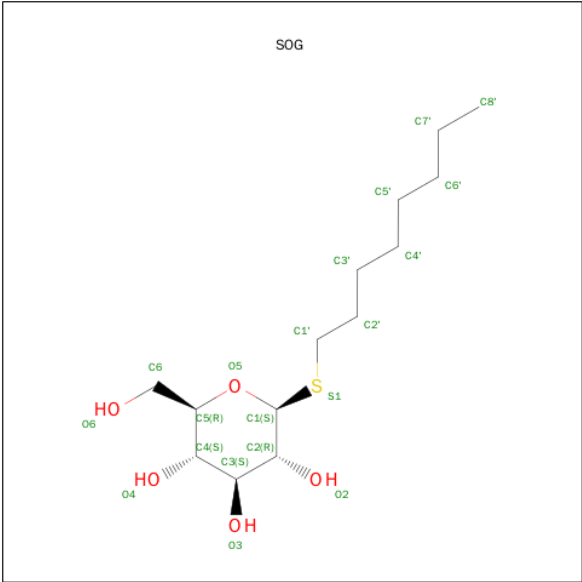
- Molecule 2 is a protein called mouse monoclonal 1gG1 Fab fragment, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	216	Total	C	N	O	S	0	0	0
			1686	1049	285	347	5			

- Molecule 3 is a protein called mouse monoclonal 1gG1 Fab fragment, kappa light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	218	Total	C	N	O	S	0	0	0
			1651	1044	278	321	8			

- Molecule 4 is N-OCTANE (three-letter code: SOG) (formula: C₁₄H₂₈O₅S).

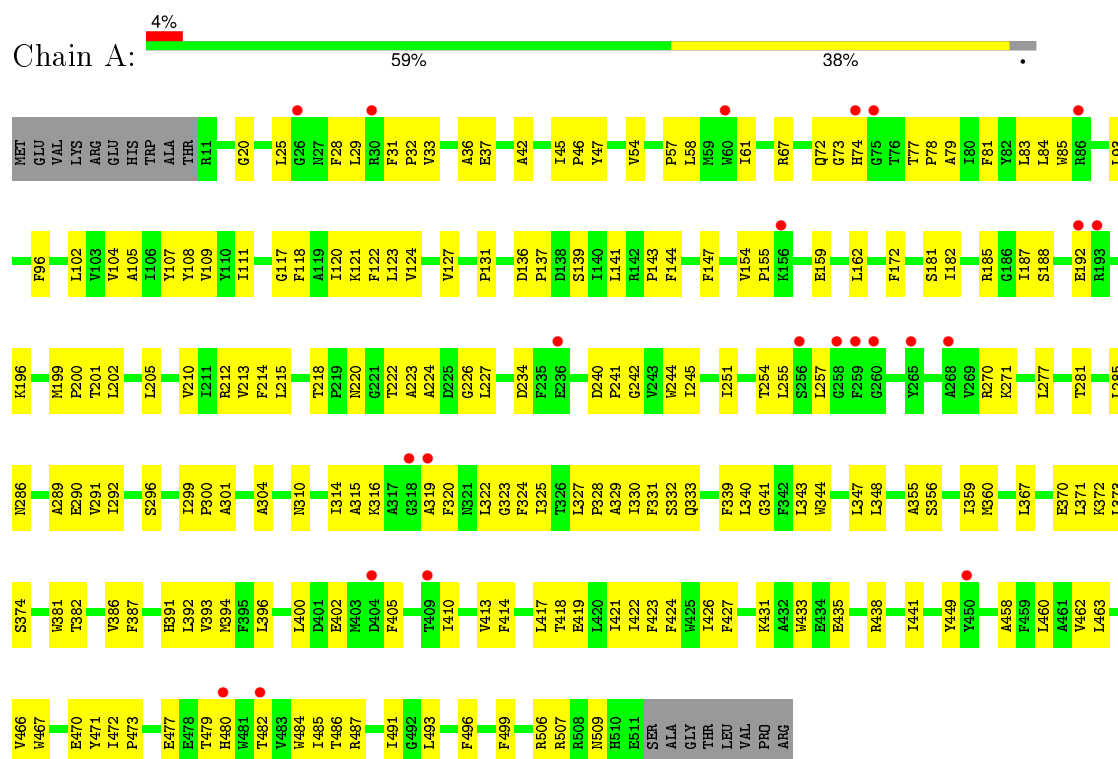


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C 8 8	0	0

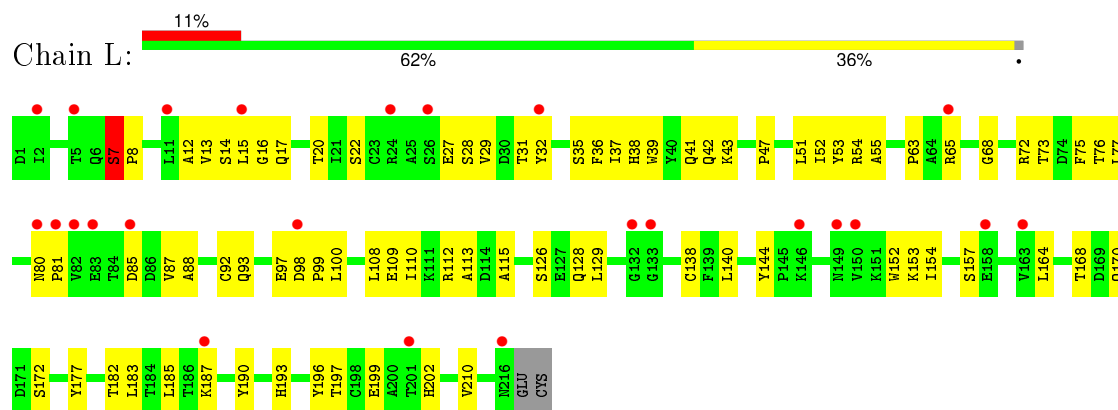
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

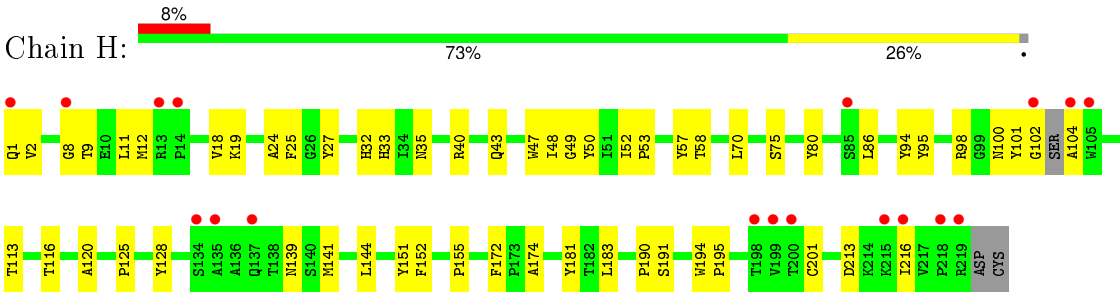
- Molecule 1: Leucine transporter LeuT



- Molecule 2: mouse monoclonal 1gG1 Fab fragment, heavy chain



- Molecule 3: mouse monoclonal 1gG1 Fab fragment, kappa light chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	128.84Å 169.80Å 130.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.58 – 3.22 35.58 – 3.22	Depositor EDS
% Data completeness (in resolution range)	90.4 (35.58-3.22) 99.0 (35.58-3.22)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 3.25Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.257 , 0.300 0.262 , 0.301	Depositor DCC
R_{free} test set	1203 reflections (5.45%)	DCC
Wilson B-factor (Å ²)	105.0	Xtriage
Anisotropy	0.612	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 65.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 23325 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7200	wwPDB-VP
Average B, all atoms (Å ²)	134.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: SOG

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	0.33	0/3976	0.64	0/5435
2	L	0.25	0/1724	0.45	0/2347
3	H	0.25	0/1694	0.47	0/2315
All	All	0.30	0/7394	0.57	0/10097

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
2	L	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	L	7	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	3855	0	3815	164	0
2	L	1686	0	1613	66	0
3	H	1651	0	1609	50	0
4	A	8	0	15	0	0
All	All	7200	0	7052	269	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 19.

All (269) 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:L:7:SER:HB3	2:L:8:PRO:CD	2.07	0.85
1:A:393:VAL:HG13	1:A:400:LEU:HB2	1.61	0.81
1:A:155:PRO:HD3	1:A:162:LEU:HD23	1.63	0.80
3:H:32:HIS:CG	3:H:98:ARG:HH21	1.99	0.80
1:A:42:ALA:HB2	1:A:234:ASP:HB3	1.65	0.77
2:L:7:SER:HB3	2:L:8:PRO:HD2	1.71	0.71
1:A:301:ALA:CB	1:A:330:ILE:HD11	2.21	0.71
2:L:37:ILE:HA	2:L:93:GLN:O	1.92	0.70
3:H:8:GLY:O	3:H:113:THR:HG23	1.93	0.69
2:L:100:LEU:HD11	3:H:47:TRP:HB2	1.75	0.69
3:H:58:THR:HG23	3:H:70:LEU:HD12	1.75	0.69
3:H:33:HIS:CD2	3:H:52:ILE:HG12	2.27	0.69
1:A:331:PHE:CE2	1:A:341:GLY:HA2	2.29	0.68
2:L:199:GLU:HG2	2:L:210:VAL:HG22	1.75	0.68
1:A:20:GLY:HA3	1:A:199:MET:HB3	1.76	0.67
3:H:32:HIS:CD2	3:H:98:ARG:HH21	2.14	0.65
1:A:332:SER:O	1:A:333:GLN:HB2	1.95	0.65
1:A:36:ALA:HB3	1:A:315:ALA:CB	2.27	0.65
1:A:188:SER:HB3	3:H:75:SER:OG	1.96	0.65
3:H:18:VAL:HG13	3:H:86:LEU:HD11	1.78	0.65
2:L:41:GLN:HB2	2:L:51:LEU:HD11	1.80	0.64
1:A:74:HIS:HE2	1:A:83:LEU:HD22	1.61	0.64
1:A:391:HIS:HA	1:A:394:MET:HG2	1.79	0.64
1:A:74:HIS:NE2	1:A:83:LEU:HD22	2.13	0.63
1:A:299:ILE:HB	1:A:300:PRO:HD3	1.81	0.63
1:A:136:ASP:HB3	1:A:139:SER:HB3	1.80	0.63
1:A:77:THR:N	1:A:78:PRO:HD2	2.14	0.63
1:A:493:LEU:HD23	1:A:493:LEU:O	1.99	0.62
2:L:183:LEU:HG	2:L:185:LEU:HD11	1.82	0.62
1:A:144:PHE:CZ	1:A:329:ALA:HA	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:458:ALA:O	1:A:462:VAL:HG23	2.00	0.61
1:A:367:LEU:HB3	1:A:373:LEU:HD12	1.81	0.61
3:H:12:MET:HG3	3:H:18:VAL:CG1	2.30	0.61
1:A:72:GLN:HG3	1:A:83:LEU:HD21	1.82	0.60
1:A:74:HIS:HB3	1:A:79:ALA:HB3	1.84	0.60
2:L:13:VAL:HG12	2:L:108:LEU:HD11	1.81	0.60
1:A:85:TRP:HB2	1:A:426:ILE:HG21	1.83	0.60
1:A:33:VAL:HG11	1:A:319:ALA:HA	1.83	0.60
1:A:58:LEU:HB2	1:A:257:LEU:HD21	1.83	0.60
1:A:58:LEU:HB3	1:A:257:LEU:HD11	1.83	0.59
3:H:11:LEU:HD12	3:H:116:THR:O	2.02	0.59
1:A:47:TYR:OH	1:A:290:GLU:OE2	2.20	0.59
2:L:42:GLN:O	2:L:88:ALA:HB1	2.02	0.59
3:H:12:MET:HG3	3:H:18:VAL:HG11	1.84	0.59
3:H:40:ARG:HB2	3:H:43:GLN:HB2	1.85	0.58
1:A:244:TRP:HB3	1:A:463:LEU:HD11	1.85	0.58
2:L:153:LYS:HB2	2:L:197:THR:HB	1.85	0.57
3:H:174:ALA:HB2	3:H:183:LEU:HD23	1.86	0.57
1:A:182:ILE:O	1:A:185:ARG:O	2.21	0.57
3:H:40:ARG:HD2	3:H:43:GLN:HG3	1.86	0.57
3:H:52:ILE:HD12	3:H:57:TYR:HD2	1.70	0.57
1:A:255:LEU:C	1:A:257:LEU:H	2.06	0.57
1:A:172:PHE:CD2	1:A:394:MET:HE1	2.39	0.57
1:A:96:PHE:CE2	1:A:414:PHE:HE2	2.22	0.57
2:L:97:GLU:O	2:L:99:PRO:O	2.23	0.57
1:A:67:ARG:NH2	1:A:435:GLU:O	2.38	0.57
1:A:136:ASP:HB3	1:A:139:SER:CB	2.34	0.56
1:A:245:ILE:HD11	1:A:413:VAL:HG11	1.87	0.56
1:A:423:PHE:O	1:A:427:PHE:HB3	2.05	0.56
1:A:74:HIS:HB3	1:A:79:ALA:CB	2.36	0.56
1:A:137:PRO:HD3	1:A:220:ASN:OD1	2.05	0.56
2:L:37:ILE:HG23	2:L:55:ALA:HB2	1.87	0.56
1:A:506:ARG:HD2	1:A:509:ASN:HD21	1.71	0.56
1:A:242:GLY:O	1:A:245:ILE:HG22	2.06	0.56
3:H:24:ALA:HB1	3:H:27:TYR:CE2	2.41	0.56
2:L:93:GLN:OE1	2:L:100:LEU:HD13	2.06	0.55
1:A:227:LEU:HD12	1:A:300:PRO:CG	2.37	0.55
3:H:141:MET:HG2	3:H:190:PRO:HA	1.87	0.55
1:A:111:ILE:HD11	1:A:400:LEU:HD11	1.89	0.55
2:L:72:ARG:HG3	2:L:73:THR:H	1.72	0.55
3:H:194:TRP:CD1	3:H:195:PRO:HA	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:174:ALA:HA	3:H:183:LEU:HB3	1.87	0.55
1:A:122:PHE:CD1	1:A:127:VAL:HG11	2.42	0.55
1:A:479:THR:HG22	1:A:484:TRP:HE1	1.71	0.55
1:A:323:GLY:O	1:A:327:LEU:HB2	2.06	0.55
1:A:370:GLU:C	1:A:372:LYS:H	2.10	0.55
2:L:27:GLU:HG2	2:L:28:SER:N	2.22	0.54
1:A:240:ASP:OD1	1:A:471:TYR:OH	2.26	0.54
2:L:98:ASP:OD1	2:L:99:PRO:HA	2.07	0.54
2:L:20:THR:HG23	2:L:76:THR:HG23	1.90	0.54
3:H:47:TRP:CZ2	3:H:49:GLY:HA2	2.43	0.54
1:A:433:TRP:CZ3	1:A:449:TYR:HB2	2.42	0.54
3:H:120:ALA:HB3	3:H:152:PHE:CE1	2.43	0.53
3:H:201:CYS:O	3:H:213:ASP:HA	2.09	0.53
1:A:421:ILE:HG13	1:A:422:ILE:N	2.24	0.53
2:L:29:VAL:O	2:L:29:VAL:HG12	2.08	0.53
1:A:487:ARG:O	1:A:491:ILE:HG12	2.07	0.53
1:A:84:LEU:HD23	1:A:423:PHE:HA	1.90	0.53
1:A:244:TRP:CB	1:A:463:LEU:HD11	2.39	0.53
1:A:290:GLU:HG3	1:A:291:VAL:N	2.24	0.52
3:H:32:HIS:O	3:H:53:PRO:HG3	2.09	0.52
1:A:201:THR:HG22	1:A:347:LEU:HD11	1.90	0.52
1:A:201:THR:CG2	1:A:347:LEU:HD11	2.40	0.52
2:L:190:TYR:HA	2:L:196:TYR:OH	2.09	0.52
1:A:105:ALA:O	1:A:109:VAL:HG23	2.10	0.52
2:L:15:LEU:HD23	2:L:16:GLY:N	2.24	0.52
2:L:39:TRP:CE2	2:L:77:LEU:HB2	2.44	0.52
1:A:28:PHE:O	1:A:32:PRO:HG2	2.10	0.52
1:A:120:ILE:O	1:A:124:VAL:HG23	2.10	0.51
1:A:81:PHE:HZ	1:A:419:GLU:HG2	1.75	0.51
1:A:141:LEU:CD1	1:A:314:ILE:HD11	2.41	0.51
2:L:12:ALA:HA	2:L:109:GLU:O	2.11	0.51
1:A:485:ILE:HG22	1:A:486:THR:N	2.26	0.51
2:L:112:ARG:HG2	2:L:113:ALA:N	2.24	0.51
1:A:393:VAL:HG13	1:A:400:LEU:CB	2.36	0.50
2:L:144:TYR:O	2:L:202:HIS:HE1	1.94	0.50
1:A:324:PHE:O	1:A:328:PRO:HG2	2.12	0.50
3:H:24:ALA:O	3:H:25:PHE:HD1	1.94	0.50
1:A:104:VAL:O	1:A:108:TYR:HB2	2.11	0.50
1:A:67:ARG:HD2	1:A:271:LYS:HA	1.93	0.50
1:A:155:PRO:HD3	1:A:162:LEU:CD2	2.38	0.50
1:A:32:PRO:HB2	1:A:322:LEU:HD23	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:52:ILE:HD13	2:L:68:GLY:N	2.27	0.50
1:A:427:PHE:O	1:A:431:LYS:HD2	2.12	0.50
2:L:80:ASN:OD1	2:L:81:PRO:HA	2.12	0.49
1:A:331:PHE:HZ	1:A:344:TRP:CD1	2.30	0.49
2:L:72:ARG:HG3	2:L:73:THR:N	2.26	0.49
1:A:154:VAL:HG13	1:A:155:PRO:HD2	1.93	0.49
1:A:42:ALA:HB2	1:A:234:ASP:CB	2.37	0.49
2:L:85:ASP:HA	2:L:172:SER:HB2	1.94	0.49
1:A:356:SER:HA	1:A:359:ILE:HG12	1.93	0.49
2:L:14:SER:HB2	2:L:17:GLN:OE1	2.12	0.49
2:L:87:VAL:CG2	2:L:110:ILE:HG12	2.42	0.49
3:H:24:ALA:HB1	3:H:27:TYR:CZ	2.47	0.49
1:A:372:LYS:HE2	1:A:507:ARG:HD3	1.94	0.49
2:L:39:TRP:CZ3	2:L:92:CYS:HB3	2.48	0.49
3:H:32:HIS:CD2	3:H:98:ARG:NH2	2.80	0.49
1:A:485:ILE:CG2	1:A:486:THR:N	2.76	0.49
1:A:320:PHE:O	1:A:324:PHE:HD2	1.96	0.49
2:L:36:PHE:O	2:L:38:HIS:CE1	2.66	0.49
1:A:292:ILE:O	1:A:296:SER:HB3	2.12	0.49
1:A:255:LEU:C	1:A:257:LEU:N	2.67	0.48
1:A:73:GLY:HA2	3:H:101:TYR:CD2	2.48	0.48
2:L:55:ALA:O	2:L:68:GLY:HA3	2.13	0.48
2:L:29:VAL:HG12	2:L:36:PHE:H	1.78	0.48
1:A:370:GLU:O	1:A:372:LYS:N	2.46	0.48
1:A:241:PRO:HB2	1:A:471:TYR:HE2	1.78	0.48
1:A:32:PRO:CB	1:A:322:LEU:HD23	2.43	0.48
1:A:496:PHE:O	1:A:499:PHE:HB3	2.14	0.48
2:L:37:ILE:CG2	2:L:55:ALA:HB2	2.44	0.47
1:A:205:LEU:HD11	1:A:340:LEU:HD11	1.97	0.47
1:A:25:LEU:CD2	1:A:348:LEU:HG	2.44	0.47
2:L:153:LYS:HA	2:L:157:SER:O	2.15	0.47
1:A:29:LEU:HD22	1:A:324:PHE:CE2	2.49	0.47
3:H:151:TYR:CE2	3:H:181:TYR:HB2	2.49	0.47
1:A:199:MET:HB2	1:A:200:PRO:HD3	1.96	0.47
2:L:47:PRO:HB3	3:H:95:TYR:CE1	2.50	0.47
1:A:210:VAL:O	1:A:213:VAL:HG12	2.15	0.47
1:A:31:PHE:HB3	1:A:32:PRO:HD3	1.97	0.47
1:A:36:ALA:CB	1:A:315:ALA:CB	2.94	0.46
2:L:37:ILE:HG21	2:L:75:PHE:CG	2.49	0.46
1:A:155:PRO:CD	1:A:162:LEU:HD23	2.40	0.46
1:A:58:LEU:C	1:A:257:LEU:HD11	2.36	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:PHE:CE2	1:A:394:MET:HE1	2.51	0.46
1:A:205:LEU:CD1	1:A:340:LEU:HD11	2.46	0.46
2:L:128:GLN:HA	3:H:128:TYR:CE1	2.51	0.46
1:A:254:THR:HG21	1:A:290:GLU:OE2	2.16	0.46
3:H:48:ILE:HD11	3:H:94:TYR:HD2	1.80	0.46
2:L:27:GLU:HG2	2:L:28:SER:H	1.80	0.45
3:H:50:TYR:O	3:H:58:THR:HG23	2.16	0.45
2:L:112:ARG:HH21	2:L:115:ALA:HB2	1.81	0.45
1:A:121:LYS:HD3	1:A:147:PHE:CZ	2.51	0.45
1:A:227:LEU:HD12	1:A:300:PRO:HG3	1.98	0.45
3:H:144:LEU:HG	3:H:216:ILE:HG21	1.98	0.45
3:H:9:THR:HG21	3:H:155:PRO:HD2	1.98	0.45
2:L:140:LEU:HD12	2:L:140:LEU:N	2.32	0.45
1:A:270:ARG:HD3	3:H:57:TYR:CE2	2.52	0.45
1:A:410:ILE:O	1:A:413:VAL:HG12	2.16	0.45
1:A:192:GLU:HA	1:A:192:GLU:OE1	2.16	0.45
1:A:386:VAL:CG2	1:A:387:PHE:N	2.80	0.45
2:L:7:SER:HB2	2:L:22:SER:HB2	1.97	0.45
2:L:168:THR:HG23	3:H:172:PHE:CD2	2.52	0.45
1:A:202:LEU:C	1:A:202:LEU:HD23	2.37	0.45
2:L:154:ILE:HD11	2:L:193:HIS:CG	2.53	0.45
2:L:164:LEU:HG	2:L:182:THR:HB	1.98	0.45
1:A:400:LEU:C	1:A:400:LEU:HD23	2.38	0.44
1:A:360:MET:CE	1:A:382:THR:HG22	2.47	0.44
2:L:126:SER:HA	2:L:129:LEU:HD12	1.99	0.44
1:A:473:PRO:O	1:A:477:GLU:HG3	2.16	0.44
1:A:45:ILE:N	1:A:46:PRO:HD2	2.32	0.44
1:A:102:LEU:HD23	1:A:493:LEU:HD21	2.00	0.44
1:A:131:PRO:HD3	1:A:143:PRO:HG3	2.00	0.44
3:H:47:TRP:HZ2	3:H:50:TYR:HD2	1.65	0.44
1:A:222:THR:HG22	1:A:224:ALA:H	1.82	0.44
1:A:466:VAL:O	1:A:470:GLU:HG2	2.17	0.44
1:A:310:ASN:O	1:A:314:ILE:HG13	2.17	0.44
1:A:325:ILE:C	1:A:328:PRO:HD2	2.38	0.44
1:A:181:SER:O	1:A:185:ARG:HG2	2.17	0.44
1:A:506:ARG:O	1:A:509:ASN:OD1	2.36	0.44
2:L:36:PHE:O	2:L:38:HIS:NE2	2.50	0.44
1:A:339:PHE:CE2	1:A:343:LEU:HD11	2.53	0.44
2:L:7:SER:O	2:L:8:PRO:C	2.56	0.44
1:A:74:HIS:CD2	1:A:83:LEU:HD22	2.53	0.44
1:A:212:ARG:O	1:A:215:LEU:HG	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:GLY:CA	1:A:199:MET:HB3	2.46	0.44
2:L:43:LYS:HG2	2:L:88:ALA:HB2	1.99	0.44
2:L:187:LYS:HB3	2:L:187:LYS:NZ	2.33	0.43
1:A:285:LEU:HA	1:A:285:LEU:HD12	1.79	0.43
1:A:480:HIS:C	1:A:482:THR:H	2.22	0.43
3:H:19:LYS:HE2	3:H:80:TYR:CD1	2.53	0.43
1:A:370:GLU:C	1:A:372:LYS:N	2.71	0.43
1:A:374:SER:HA	3:H:1:GLN:CD	2.38	0.43
1:A:493:LEU:HD23	1:A:493:LEU:C	2.38	0.43
1:A:187:ILE:HG13	1:A:187:ILE:O	2.17	0.43
1:A:93:LEU:HB3	1:A:418:THR:HG21	1.99	0.43
1:A:286:ASN:O	1:A:289:ALA:HB3	2.18	0.43
1:A:277:LEU:HD23	1:A:441:ILE:HD12	2.00	0.43
2:L:37:ILE:HG21	2:L:75:PHE:CD2	2.54	0.43
2:L:99:PRO:HB3	3:H:47:TRP:HZ3	1.84	0.43
1:A:202:LEU:O	1:A:202:LEU:HD23	2.18	0.43
3:H:47:TRP:HZ2	3:H:50:TYR:CD2	2.36	0.43
1:A:37:GLU:HG2	1:A:316:LYS:HA	2.01	0.43
1:A:382:THR:O	1:A:386:VAL:HG13	2.19	0.43
2:L:53:TYR:HB2	3:H:104:ALA:HB2	2.01	0.43
1:A:506:ARG:HA	1:A:509:ASN:ND2	2.34	0.43
1:A:25:LEU:HD21	1:A:348:LEU:HG	2.01	0.43
1:A:405:PHE:HZ	1:A:467:TRP:HZ2	1.66	0.43
1:A:141:LEU:HD11	1:A:314:ILE:HD11	2.01	0.43
1:A:213:VAL:HG13	1:A:214:PHE:N	2.34	0.42
3:H:139:ASN:O	3:H:191:SER:HB3	2.19	0.42
1:A:392:LEU:O	1:A:396:LEU:HB2	2.19	0.42
1:A:118:PHE:CE2	1:A:324:PHE:HD1	2.37	0.42
2:L:126:SER:O	2:L:129:LEU:HB2	2.19	0.42
2:L:170:GLN:HB2	2:L:177:TYR:CZ	2.54	0.42
1:A:58:LEU:CB	1:A:257:LEU:HD11	2.50	0.42
1:A:141:LEU:HD13	1:A:314:ILE:HD11	2.01	0.42
1:A:356:SER:HA	1:A:359:ILE:CG1	2.49	0.42
1:A:226:GLY:HA3	1:A:300:PRO:HA	2.01	0.42
1:A:343:LEU:O	1:A:347:LEU:HG	2.20	0.42
1:A:325:ILE:O	1:A:328:PRO:HD2	2.19	0.42
3:H:125:PRO:HB3	3:H:151:TYR:HB3	2.01	0.42
1:A:54:VAL:O	1:A:58:LEU:HD13	2.20	0.42
3:H:100:ASN:C	3:H:102:GLY:H	2.22	0.42
2:L:43:LYS:HE2	2:L:85:ASP:O	2.20	0.42
3:H:2:VAL:HG22	3:H:27:TYR:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:417:LEU:HD12	1:A:460:LEU:HB3	2.02	0.42
2:L:100:LEU:CD1	3:H:47:TRP:HB2	2.48	0.42
2:L:138:CYS:HB2	2:L:152:TRP:CZ2	2.54	0.42
1:A:417:LEU:C	1:A:417:LEU:HD23	2.40	0.42
3:H:35:ASN:OD1	3:H:50:TYR:HB3	2.20	0.42
1:A:402:GLU:HA	1:A:402:GLU:OE1	2.20	0.42
2:L:87:VAL:O	2:L:88:ALA:HB2	2.20	0.41
2:L:63:PRO:HB2	2:L:65:ARG:HG3	2.01	0.41
1:A:84:LEU:HG	1:A:427:PHE:CD1	2.54	0.41
1:A:251:ILE:HA	1:A:251:ILE:HD13	1.93	0.41
1:A:155:PRO:HB2	1:A:159:GLU:HB2	2.01	0.41
1:A:57:PRO:O	1:A:61:ILE:HG12	2.20	0.41
1:A:281:THR:O	1:A:285:LEU:HB2	2.21	0.41
1:A:77:THR:N	1:A:78:PRO:CD	2.82	0.41
2:L:185:LEU:N	2:L:185:LEU:HD12	2.36	0.41
3:H:57:TYR:O	3:H:57:TYR:CG	2.74	0.41
2:L:93:GLN:HE21	2:L:93:GLN:HB3	1.66	0.41
2:L:98:ASP:HA	2:L:99:PRO:C	2.41	0.41
1:A:396:LEU:CD2	1:A:485:ILE:HG21	2.51	0.41
1:A:355:ALA:O	1:A:359:ILE:HG12	2.21	0.41
3:H:48:ILE:HD11	3:H:94:TYR:CD2	2.55	0.41
1:A:192:GLU:O	1:A:196:LYS:HG3	2.21	0.41
2:L:35:SER:O	2:L:54:ARG:HA	2.20	0.41
1:A:33:VAL:O	1:A:37:GLU:HG3	2.21	0.41
1:A:360:MET:HE3	1:A:382:THR:HG22	2.03	0.41
1:A:472:ILE:CB	1:A:473:PRO:HD3	2.51	0.41
1:A:104:VAL:HA	1:A:107:TYR:CE2	2.55	0.41
2:L:31:THR:O	2:L:32:TYR:HB2	2.21	0.41
1:A:290:GLU:HG3	1:A:291:VAL:H	1.85	0.40
1:A:435:GLU:CD	1:A:438:ARG:HH21	2.24	0.40
1:A:223:ALA:HA	1:A:304:ALA:HB2	2.03	0.40
1:A:120:ILE:O	1:A:123:LEU:HB3	2.22	0.40
1:A:373:LEU:HD11	1:A:381:TRP:CZ3	2.56	0.40
1:A:117:GLY:O	1:A:121:LYS:HG3	2.21	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	499/519 (96%)	479 (96%)	19 (4%)	1 (0%)	52	88
2	L	214/218 (98%)	205 (96%)	8 (4%)	1 (0%)	34	77
3	H	214/221 (97%)	203 (95%)	11 (5%)	0	100	100
All	All	927/958 (97%)	887 (96%)	38 (4%)	2 (0%)	52	88

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	371	LEU
2	L	7	SER

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	373/417 (89%)	371 (100%)	2 (0%)	92	97
2	L	194/196 (99%)	193 (100%)	1 (0%)	92	97
3	H	184/189 (97%)	184 (100%)	0	100	100
All	All	751/802 (94%)	748 (100%)	3 (0%)	93	98

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	218	THR

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Mol	Chain	Res	Type
1	A	424	PHE
2	L	7	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	361	GLN
3	H	39	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z > 2$
4	SOG	A	520	-	7,7,20	0.24	0	6,6,25	0.66	0

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
4	SOG	A	520	-	-	0/5/5/31	0/0/0/1

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](#)

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	501/519 (96%)	0.31	23 (4%) 36 24	75, 117, 166, 218	0
2	L	216/218 (99%)	0.66	24 (11%) 7 5	104, 152, 186, 227	0
3	H	218/221 (98%)	0.51	18 (8%) 14 8	99, 146, 197, 227	0
All	All	935/958 (97%)	0.44	65 (6%) 19 12	75, 132, 181, 227	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	219	ARG	5.2
3	H	215	LYS	4.9
2	L	216	ASN	4.8
2	L	201	THR	4.8
2	L	15	LEU	4.2
1	A	74	HIS	4.1
1	A	258	GLY	4.1
3	H	104	ALA	4.1
1	A	192	GLU	4.1
1	A	260	GLY	3.9
3	H	216	ILE	3.9
1	A	404	ASP	3.9
2	L	24	ARG	3.9
1	A	319	ALA	3.8
3	H	102	GLY	3.7
3	H	135	ALA	3.6
3	H	134	SER	3.6
2	L	146	LYS	3.6
1	A	236	GLU	3.6
2	L	132	GLY	3.6
2	L	163	VAL	3.4
2	L	187	LYS	3.2
2	L	5	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	256	SER	2.9
2	L	32	TYR	2.9
2	L	80	ASN	2.8
2	L	65	ARG	2.8
2	L	83	GLU	2.8
2	L	81	PRO	2.8
1	A	156	LYS	2.8
1	A	26	GLY	2.8
1	A	409	THR	2.7
1	A	86	ARG	2.7
3	H	218	PRO	2.7
1	A	193	ARG	2.7
3	H	13	ARG	2.6
1	A	75	GLY	2.6
3	H	105	TRP	2.6
2	L	149	ASN	2.6
2	L	2	ILE	2.5
2	L	11	LEU	2.5
3	H	200	THR	2.4
1	A	30	ARG	2.4
3	H	14	PRO	2.4
2	L	133	GLY	2.4
2	L	85	ASP	2.3
3	H	199	VAL	2.3
3	H	8	GLY	2.3
3	H	85	SER	2.3
3	H	1	GLN	2.2
1	A	60	TRP	2.2
2	L	26	SER	2.2
2	L	158	GLU	2.2
1	A	482	THR	2.2
2	L	150	VAL	2.2
3	H	198	THR	2.1
1	A	480	HIS	2.1
1	A	268	ALA	2.1
1	A	259	PHE	2.1
1	A	450	TYR	2.1
2	L	98	ASP	2.1
1	A	318	GLY	2.1
1	A	265	TYR	2.0
2	L	82	VAL	2.0
3	H	137	GLN	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
4	SOG	A	520	8/20	0.66	0.94	15.81	84,112,136,144	0

6.5 Other polymers [i](#)

There are no such residues in this entry.