



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:31 PM GMT

PDB ID : 4ILS
Title : Crystal structure of engineered protein. northeast structural genomics Consortium target or117
Authors : Seetharaman, J.; Lew, S.; Nivon, L.; Baker, D.; Bjelic, S.; Ciccocanti, C.; Sahdev, S.; Xiao, R.; Everett, J.K.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2012-12-31
Resolution : 2.50 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

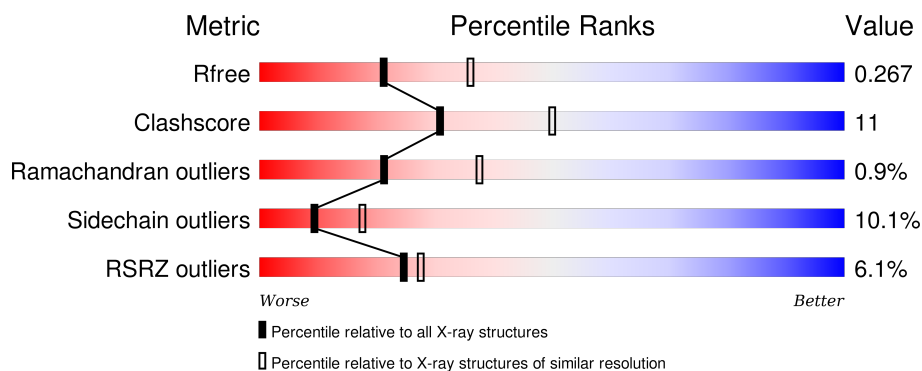
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	390	<div> <div>4%</div> <div>68% 18% • 10%</div> </div>
1	B	390	<div> <div>4%</div> <div>65% 18% 5% 12%</div> </div>
1	C	390	<div> <div>8%</div> <div>63% 21% • 13%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8504 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 Engineered protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	350	Total	C	N	O	S	0	0	0
			2786	1792	490	490	14			
1	B	344	Total	C	N	O	S	0	0	0
			2726	1749	481	481	15			
1	C	341	Total	C	N	O	S	0	0	0
			2721	1748	480	479	14			

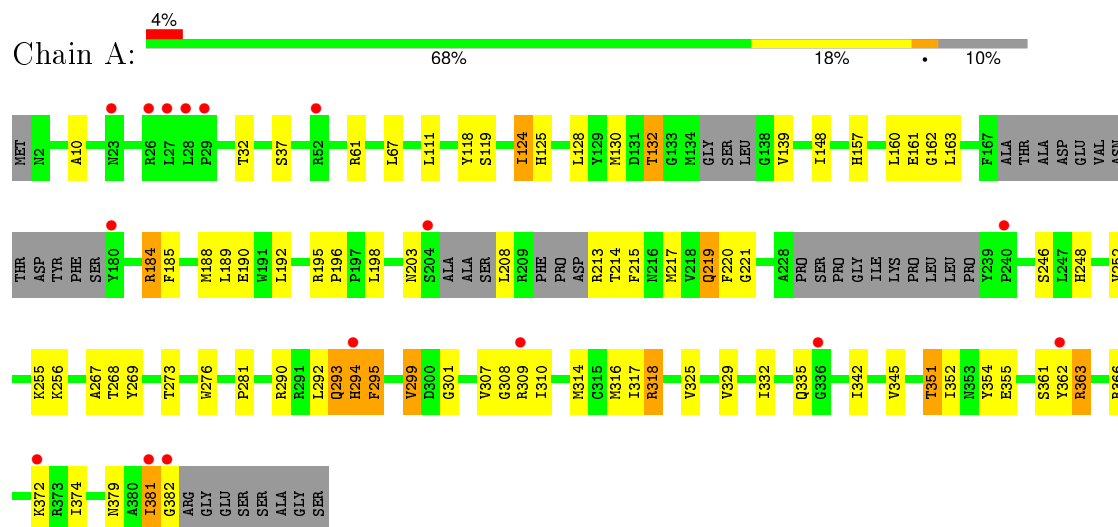
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	105	Total	O	0	0
			105	105		
2	B	88	Total	O	0	0
			88	88		
2	C	78	Total	O	0	0
			78	78		

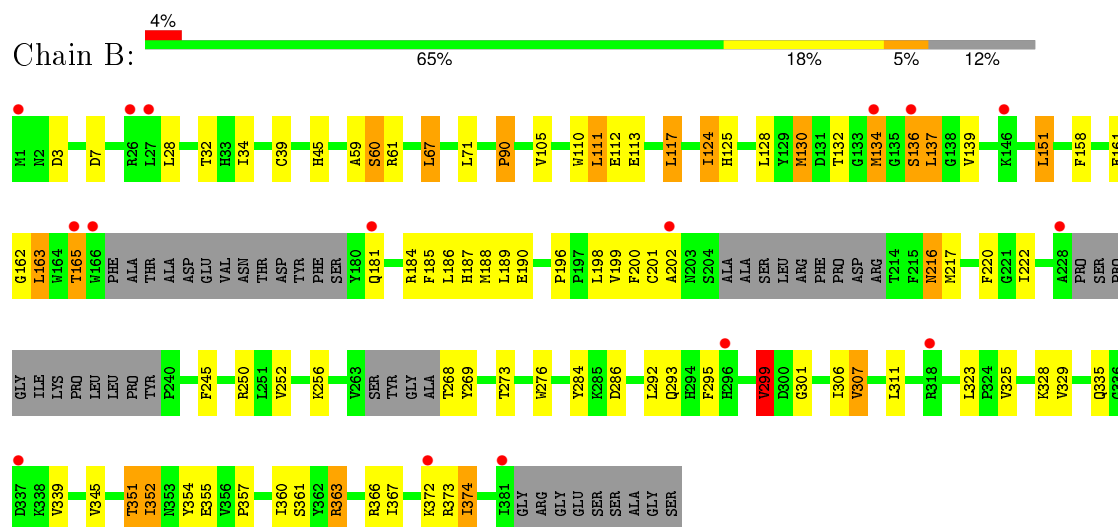
3 Residue-property plots

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: Engineered protein

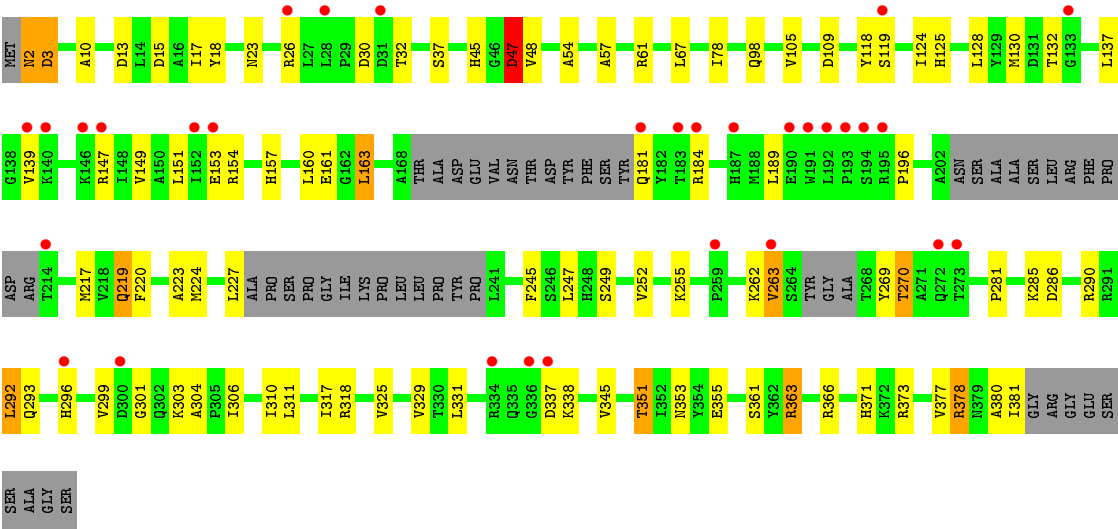


• Molecule 1: Engineered protein



• Molecule 1: Engineered protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	112.85Å 112.85Å 232.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.37 – 2.50 49.32 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.37-2.50) 99.6 (49.32-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.75 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.218 , 0.265 0.221 , 0.267	Depositor DCC
R_{free} test set	2679 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	42.8	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 52655 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8504	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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](#)

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	0.49	0/2854	0.68	1/3871 (0.0%)
1	B	0.46	0/2791	0.67	1/3783 (0.0%)
1	C	0.39	0/2786	0.63	0/3777
All	All	0.45	0/8431	0.66	2/11431 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	293	GLN	N-CA-C	-6.12	94.47	111.00
1	B	299	VAL	CB-CA-C	-5.77	100.44	111.40

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	A	2786	0	2756	59	0
1	B	2726	0	2716	74	0
1	C	2721	0	2720	53	0
2	A	105	0	0	8	0
2	B	88	0	0	3	0
2	C	78	0	0	5	0
All	All	8504	0	8192	179	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 11.

All (179) 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:B:130:MET:O	1:B:165:THR:HG22	1.42	1.16
1:B:134:MET:HB2	1:B:181:GLN:NE2	1.64	1.09
1:A:381:ILE:HD11	2:A:439:HOH:O	1.58	1.03
1:B:134:MET:CB	1:B:181:GLN:NE2	2.26	0.97
1:C:147:ARG:HG2	2:C:470:HOH:O	1.68	0.93
1:B:134:MET:HB2	1:B:181:GLN:HE22	1.31	0.92
1:A:132:THR:HG22	2:A:404:HOH:O	1.72	0.89
1:A:128:LEU:HB3	1:A:130:MET:CE	2.03	0.88
1:C:361:SER:OG	1:C:363:ARG:HG3	1.75	0.86
1:B:130:MET:O	1:B:165:THR:CG2	2.24	0.85
1:B:299:VAL:HG13	1:B:329:VAL:HG22	1.58	0.85
1:A:351:THR:HG21	1:A:355:GLU:OE2	1.79	0.83
1:B:374:ILE:HG22	2:B:429:HOH:O	1.78	0.82
1:B:351:THR:HG22	1:C:290:ARG:HH11	1.44	0.79
1:A:190:GLU:OE2	1:B:187:HIS:CE1	2.36	0.79
1:B:124:ILE:CD1	1:B:158:PHE:CD1	2.67	0.77
1:A:128:LEU:HB3	1:A:130:MET:HE2	1.66	0.77
1:B:130:MET:HB2	1:B:165:THR:HG23	1.68	0.75
1:A:190:GLU:OE2	1:B:187:HIS:HE1	1.70	0.74
1:B:199:VAL:H	1:B:216:ASN:HD21	1.37	0.73
1:A:379:ASN:OD1	1:A:381:ILE:HG13	1.88	0.73
1:C:351:THR:HG21	1:C:355:GLU:OE2	1.88	0.73
1:A:293:GLN:O	1:A:294:HIS:HB2	1.89	0.72
1:B:199:VAL:H	1:B:216:ASN:ND2	1.88	0.72
1:B:132:THR:HG23	1:B:165:THR:HG21	1.73	0.71
1:B:286:ASP:OD1	1:B:363:ARG:HD2	1.90	0.71
1:C:48:VAL:HG23	1:C:78:ILE:CD1	2.21	0.70
1:A:361:SER:OG	1:A:363:ARG:HG3	1.91	0.70
1:C:48:VAL:HG23	1:C:78:ILE:HD13	1.73	0.69
1:C:3:ASP:HB3	2:C:420:HOH:O	1.93	0.68
1:B:134:MET:HA	1:B:181:GLN:HE21	1.59	0.68
1:C:2:ASN:HB2	1:C:377:VAL:O	1.94	0.68
1:B:301:GLY:H	1:B:335:GLN:HE21	1.44	0.66
1:B:128:LEU:HB3	1:B:130:MET:CE	2.26	0.65
1:C:286:ASP:OD1	1:C:363:ARG:HD3	1.97	0.64
1:A:125:HIS:HD2	1:A:161:GLU:OE2	1.79	0.64
1:A:299:VAL:HG13	1:A:329:VAL:HG22	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:HIS:CD2	1:A:161:GLU:OE2	2.51	0.64
1:C:54:ALA:HB2	1:C:224:MET:CE	2.29	0.63
1:B:128:LEU:HB3	1:B:130:MET:HE2	1.81	0.63
1:B:351:THR:HG21	1:B:355:GLU:OE2	1.99	0.63
1:A:290:ARG:HA	1:A:310:ILE:HB	1.81	0.63
1:C:363:ARG:HB3	1:C:381:ILE:HD11	1.81	0.62
1:B:136:SER:HB2	1:B:137:LEU:HD13	1.81	0.61
1:C:132:THR:O	1:C:181:GLN:HG2	2.01	0.61
1:B:124:ILE:HD11	1:B:158:PHE:CD1	2.36	0.61
1:C:125:HIS:HD2	1:C:161:GLU:OE2	1.83	0.60
1:A:163:LEU:HD13	1:A:192:LEU:HD11	1.83	0.60
1:A:37:SER:HB2	1:A:219:GLN:HE22	1.66	0.60
1:A:301:GLY:H	1:A:335:GLN:HE21	1.49	0.59
1:A:61:ARG:NH2	1:A:161:GLU:OE1	2.35	0.59
1:C:54:ALA:HB2	1:C:224:MET:HE3	1.83	0.59
1:B:130:MET:CB	1:B:165:THR:HG23	2.33	0.59
1:B:3:ASP:HB3	2:B:486:HOH:O	2.03	0.59
1:B:134:MET:CA	1:B:181:GLN:HE21	2.16	0.58
1:A:128:LEU:HD21	1:A:148:ILE:HG21	1.85	0.58
1:A:352:ILE:HD13	1:A:354:TYR:HB2	1.84	0.58
1:C:361:SER:HG	1:C:363:ARG:HG3	1.70	0.57
1:A:292:LEU:HD21	1:A:345:VAL:HG13	1.85	0.57
1:B:201:CYS:O	1:B:202:ALA:HB3	2.04	0.57
1:A:214:THR:HG22	2:A:412:HOH:O	2.04	0.57
1:A:293:GLN:HG3	1:A:308:GLY:O	2.04	0.57
1:C:10:ALA:HB3	1:C:366:ARG:HD2	1.85	0.57
1:C:61:ARG:HD2	2:C:462:HOH:O	2.04	0.57
1:B:39:CYS:SG	1:C:285:LYS:NZ	2.78	0.57
1:B:222:ILE:HG12	1:B:354:TYR:CZ	2.40	0.56
1:A:293:GLN:O	1:A:294:HIS:CB	2.54	0.56
1:B:90:PRO:HD3	1:B:110:TRP:NE1	2.21	0.55
1:B:256:LYS:HD3	1:B:276:TRP:CH2	2.41	0.55
1:A:374:ILE:HG22	2:A:465:HOH:O	2.05	0.55
1:B:269:TYR:CE1	1:B:307:VAL:HG22	2.41	0.55
1:B:134:MET:CB	1:B:181:GLN:HE21	2.14	0.54
1:A:221:GLY:N	2:A:444:HOH:O	2.39	0.54
1:B:124:ILE:HD13	1:B:158:PHE:CD1	2.40	0.54
1:A:318:ARG:HD2	2:A:496:HOH:O	2.07	0.54
1:C:227:LEU:HD23	1:C:353:ASN:HD21	1.71	0.54
1:A:290:ARG:NH2	1:A:309:ARG:HD2	2.22	0.54
1:B:301:GLY:H	1:B:335:GLN:NE2	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:367:ILE:HD13	1:B:367:ILE:N	2.23	0.54
1:A:362:TYR:H	1:A:362:TYR:HD2	1.54	0.54
1:B:216:ASN:H	1:B:216:ASN:HD22	1.56	0.54
1:A:128:LEU:HD13	1:A:160:LEU:HD13	1.89	0.54
1:A:281:PRO:HA	1:A:314:MET:HG2	1.90	0.54
1:B:130:MET:CB	1:B:165:THR:CG2	2.86	0.54
1:A:381:ILE:O	1:A:382:GLY:C	2.45	0.53
1:C:304:ALA:HB1	1:C:318:ARG:O	2.08	0.53
1:B:351:THR:HG22	1:C:290:ARG:NH1	2.19	0.53
1:B:163:LEU:HB2	1:B:196:PRO:HG2	1.90	0.53
1:B:61:ARG:HD3	1:B:217:MET:SD	2.49	0.53
1:A:269:TYR:CD1	1:A:307:VAL:HG13	2.44	0.53
1:B:34:ILE:HB	1:B:59:ALA:HA	1.91	0.52
1:C:125:HIS:CD2	1:C:161:GLU:OE2	2.63	0.52
1:C:128:LEU:HB3	1:C:130:MET:CE	2.39	0.52
1:C:154:ARG:HD2	2:C:416:HOH:O	2.10	0.52
1:C:292:LEU:HD21	1:C:345:VAL:HG13	1.90	0.52
1:C:303:LYS:HG3	1:C:338:LYS:HE3	1.93	0.51
1:B:295:PHE:HB3	1:B:306:ILE:HD12	1.92	0.51
1:B:198:LEU:HA	1:B:216:ASN:HD21	1.77	0.50
1:A:381:ILE:CD1	2:A:439:HOH:O	2.37	0.50
1:A:316:MET:HB3	2:A:464:HOH:O	2.11	0.50
1:B:184:ARG:O	1:B:188:MET:HG3	2.11	0.49
1:C:249:SER:OG	1:C:281:PRO:HD2	2.13	0.49
1:C:247:LEU:HG	1:C:331:LEU:HD23	1.94	0.49
1:A:214:THR:HA	1:A:215:PHE:HB2	1.95	0.48
1:A:379:ASN:OD1	1:A:381:ILE:CG1	2.60	0.48
1:B:45:HIS:NE2	1:B:245:PHE:HB2	2.29	0.48
1:B:165:THR:OG1	1:B:185:PHE:CE1	2.65	0.48
1:B:124:ILE:HD11	1:B:158:PHE:CE1	2.48	0.48
1:A:362:TYR:CD2	1:A:362:TYR:N	2.80	0.47
1:B:128:LEU:HB3	1:B:130:MET:HE1	1.95	0.47
1:C:47:ASP:HB3	1:C:48:VAL:H	1.41	0.47
1:C:2:ASN:HB3	1:C:3:ASP:H	1.53	0.47
1:B:60:SER:HB2	1:B:61:ARG:HG3	1.96	0.46
1:C:119:SER:O	1:C:157:HIS:CD2	2.68	0.46
1:B:7:ASP:HB3	1:B:250:ARG:HB2	1.97	0.46
1:A:190:GLU:CD	1:B:187:HIS:CE1	2.89	0.46
1:A:61:ARG:HD3	1:A:217:MET:SD	2.55	0.46
1:A:119:SER:O	1:A:157:HIS:HD2	1.98	0.46
1:C:18:TYR:CD2	1:C:57:ALA:HB2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:ASP:OD1	1:C:15:ASP:HB2	2.15	0.46
1:C:54:ALA:HB2	1:C:224:MET:HE1	1.97	0.46
1:C:263:VAL:HG22	1:C:269:TYR:HB3	1.97	0.46
1:C:223:ALA:HA	1:C:227:LEU:HD12	1.98	0.45
1:A:246:SER:HB3	1:A:248:HIS:NE2	2.32	0.45
1:A:301:GLY:H	1:A:335:GLN:NE2	2.14	0.45
1:B:128:LEU:HB2	1:B:163:LEU:HD22	1.98	0.45
1:C:163:LEU:HB2	1:C:196:PRO:HG3	1.98	0.45
1:A:267:ALA:HB2	1:A:309:ARG:HG2	1.98	0.45
1:B:124:ILE:CD1	1:B:158:PHE:HD1	2.23	0.44
1:C:371:HIS:C	1:C:373:ARG:H	2.19	0.44
1:B:113:GLU:O	1:B:117:LEU:HG	2.18	0.44
1:B:352:ILE:H	1:B:352:ILE:HG13	1.43	0.44
1:B:162:GLY:HA2	1:B:198:LEU:O	2.17	0.44
1:A:332:ILE:HG21	1:A:342:ILE:HD13	1.99	0.44
1:C:61:ARG:HD3	1:C:217:MET:SD	2.58	0.44
1:A:213:ARG:C	1:A:214:THR:HG23	2.37	0.44
1:C:378:ARG:HH11	1:C:380:ALA:HA	1.82	0.44
1:A:162:GLY:HA2	1:A:198:LEU:O	2.18	0.44
1:B:130:MET:HE3	1:B:163:LEU:CD1	2.48	0.43
1:C:45:HIS:NE2	1:C:245:PHE:HB2	2.33	0.43
1:A:10:ALA:HB3	1:A:366:ARG:HD2	2.00	0.43
1:C:23:ASN:O	1:C:26:ARG:HB3	2.18	0.43
1:B:28:LEU:HD12	1:B:34:ILE:HD11	2.00	0.43
1:A:267:ALA:HB2	1:A:309:ARG:CG	2.48	0.43
1:C:296:HIS:HB3	2:C:464:HOH:O	2.18	0.43
1:B:125:HIS:HD2	1:B:161:GLU:OE2	2.01	0.43
1:B:222:ILE:H	1:B:222:ILE:HG13	1.64	0.43
1:B:67:LEU:HD22	1:B:71:LEU:HG	2.01	0.43
1:A:290:ARG:HH21	1:A:309:ARG:HD2	1.83	0.43
1:C:262:LYS:HG2	1:C:270:THR:HG23	2.00	0.43
1:A:203:ASN:O	1:A:208:LEU:N	2.52	0.42
1:A:267:ALA:CB	1:A:309:ARG:HG2	2.49	0.42
1:B:111:LEU:HA	1:B:111:LEU:HD23	1.94	0.42
1:A:128:LEU:HD23	1:A:130:MET:HE3	2.00	0.42
1:B:124:ILE:CD1	1:B:158:PHE:CE1	3.02	0.42
1:A:163:LEU:HD23	1:A:185:PHE:CZ	2.54	0.42
1:B:130:MET:HE1	1:B:163:LEU:CD2	2.49	0.42
1:C:128:LEU:HB3	1:C:130:MET:HE3	2.02	0.42
1:C:378:ARG:NH1	1:C:380:ALA:HA	2.34	0.42
1:A:256:LYS:HD3	1:A:276:TRP:CZ2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:MET:HE1	1:B:163:LEU:HD21	2.02	0.41
1:B:134:MET:CA	1:B:181:GLN:NE2	2.77	0.41
1:B:293:GLN:HB3	2:B:409:HOH:O	2.19	0.41
1:C:128:LEU:HD13	1:C:160:LEU:CD1	2.50	0.41
1:B:292:LEU:HD21	1:B:345:VAL:HG13	2.02	0.41
1:A:195:ARG:HA	1:A:196:PRO:HD3	1.93	0.41
1:C:306:ILE:HA	1:C:317:ILE:HG22	2.02	0.41
1:C:299:VAL:HG13	1:C:329:VAL:HG22	2.02	0.41
1:B:200:PHE:HB3	1:B:217:MET:HB3	2.03	0.41
1:C:118:TYR:CD2	1:C:124:ILE:HD11	2.56	0.40
1:A:184:ARG:O	1:A:188:MET:HG3	2.20	0.40
1:A:118:TYR:CD2	1:A:124:ILE:HD11	2.56	0.40
1:B:28:LEU:HD12	1:B:34:ILE:CD1	2.51	0.40
1:C:37:SER:HB2	1:C:219:GLN:OE1	2.21	0.40
1:B:284:TYR:CD2	1:C:355:GLU:HG3	2.56	0.40
1:A:294:HIS:O	1:A:295:PHE:O	2.39	0.40
1:B:112:GLU:HG3	1:B:151:LEU:HD21	2.02	0.40
1:C:149:VAL:O	1:C:153:GLU:HG3	2.21	0.40
1:B:357:PRO:HA	1:B:360:ILE:HD12	2.02	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	338/390 (87%)	327 (97%)	7 (2%)	4 (1%)	16	29
1	B	334/390 (86%)	318 (95%)	16 (5%)	0	100	100
1	C	331/390 (85%)	313 (95%)	13 (4%)	5 (2%)	13	22
All	All	1003/1170 (86%)	958 (96%)	36 (4%)	9 (1%)	21	37

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	47	ASP
1	A	294	HIS
1	A	268	THR
1	C	263	VAL
1	C	301	GLY
1	C	293	GLN
1	A	295	PHE
1	A	372	LYS
1	C	3	ASP

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	289/325 (89%)	269 (93%)	20 (7%)	19	35
1	B	286/325 (88%)	247 (86%)	39 (14%)	5	8
1	C	287/325 (88%)	259 (90%)	28 (10%)	10	19
All	All	862/975 (88%)	775 (90%)	87 (10%)	9	17

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	THR
1	A	67	LEU
1	A	111	LEU
1	A	124	ILE
1	A	132	THR
1	A	139	VAL
1	A	184	ARG
1	A	189	LEU
1	A	219	GLN
1	A	220	PHE
1	A	252	VAL
1	A	255	LYS
1	A	273	THR
1	A	299	VAL

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Mol	Chain	Res	Type
1	A	317	ILE
1	A	318	ARG
1	A	325	VAL
1	A	351	THR
1	A	363	ARG
1	A	381	ILE
1	B	32	THR
1	B	60	SER
1	B	67	LEU
1	B	90	PRO
1	B	105	VAL
1	B	111	LEU
1	B	117	LEU
1	B	124	ILE
1	B	130	MET
1	B	134	MET
1	B	136	SER
1	B	137	LEU
1	B	139	VAL
1	B	151	LEU
1	B	163	LEU
1	B	165	THR
1	B	186	LEU
1	B	189	LEU
1	B	190	GLU
1	B	216	ASN
1	B	220	PHE
1	B	252	VAL
1	B	268	THR
1	B	273	THR
1	B	299	VAL
1	B	307	VAL
1	B	311	LEU
1	B	323	LEU
1	B	325	VAL
1	B	328	LYS
1	B	339	VAL
1	B	351	THR
1	B	352	ILE
1	B	361	SER
1	B	363	ARG
1	B	366	ARG

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Mol	Chain	Res	Type
1	B	372	LYS
1	B	373	ARG
1	B	374	ILE
1	C	2	ASN
1	C	17	ILE
1	C	30	ASP
1	C	32	THR
1	C	47	ASP
1	C	67	LEU
1	C	98	GLN
1	C	105	VAL
1	C	109	ASP
1	C	137	LEU
1	C	139	VAL
1	C	151	LEU
1	C	163	LEU
1	C	184	ARG
1	C	189	LEU
1	C	219	GLN
1	C	220	PHE
1	C	252	VAL
1	C	255	LYS
1	C	270	THR
1	C	292	LEU
1	C	310	ILE
1	C	311	LEU
1	C	325	VAL
1	C	337	ASP
1	C	351	THR
1	C	363	ARG
1	C	378	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	HIS
1	A	157	HIS
1	A	203	ASN
1	A	219	GLN
1	A	258	GLN
1	A	335	GLN
1	B	125	HIS

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Mol	Chain	Res	Type
1	B	157	HIS
1	B	181	GLN
1	B	187	HIS
1	B	203	ASN
1	B	216	ASN
1	B	258	GLN
1	B	335	GLN
1	C	125	HIS
1	C	157	HIS
1	C	335	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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	350/390 (89%)	0.25	16 (4%)	36 41	20, 36, 62, 75	0
1	B	344/390 (88%)	0.34	16 (4%)	35 40	22, 38, 63, 107	0
1	C	341/390 (87%)	0.50	31 (9%)	11 12	25, 46, 74, 94	0
All	All	1035/1170 (88%)	0.36	63 (6%)	25 27	20, 40, 68, 107	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	MET	6.3
1	B	337	ASP	4.5
1	B	228	ALA	3.7
1	C	272	GLN	3.6
1	C	26	ARG	3.5
1	A	180	TYR	3.5
1	B	296	HIS	3.4
1	C	190	GLU	3.4
1	C	214	THR	3.4
1	C	334	ARG	3.4
1	A	382	GLY	3.3
1	C	133	GLY	3.3
1	C	336	GLY	3.3
1	C	192	LEU	3.3
1	C	31	ASP	3.2
1	A	26	ARG	3.1
1	B	134	MET	3.1
1	B	27	LEU	3.1
1	C	296	HIS	3.0
1	B	26	ARG	3.0
1	C	187	HIS	3.0
1	C	300	ASP	3.0
1	A	204	SER	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	136	SER	2.9
1	C	153	GLU	2.9
1	A	23	ASN	2.8
1	A	27	LEU	2.7
1	A	309	ARG	2.7
1	B	381	ILE	2.7
1	A	28	LEU	2.7
1	A	52	ARG	2.7
1	B	166	TRP	2.7
1	B	146	LYS	2.7
1	A	294	HIS	2.6
1	C	181	GLN	2.6
1	C	183	THR	2.5
1	C	119	SER	2.5
1	A	240	PRO	2.5
1	A	29	PRO	2.5
1	C	193	PRO	2.4
1	B	372	LYS	2.4
1	C	184	ARG	2.4
1	C	191	TRP	2.4
1	B	165	THR	2.4
1	C	195	ARG	2.4
1	A	362	TYR	2.4
1	C	146	LYS	2.4
1	C	259	PRO	2.3
1	C	147	ARG	2.3
1	A	381	ILE	2.3
1	C	273	THR	2.3
1	C	140	LYS	2.3
1	C	194	SER	2.3
1	B	318	ARG	2.3
1	C	28	LEU	2.3
1	C	337	ASP	2.3
1	C	263	VAL	2.2
1	C	139	VAL	2.2
1	C	152	ILE	2.1
1	B	181	GLN	2.1
1	B	202	ALA	2.1
1	A	372	LYS	2.1
1	A	336	GLY	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.