



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:22 AM UTC

PDB ID : 8K2Y / pdb_00008k2y
Title : Crystal structure of MucD
Authors : Kim, J.H.; Park, H.H.
Deposited on : 2023-07-14
Resolution : 2.70 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

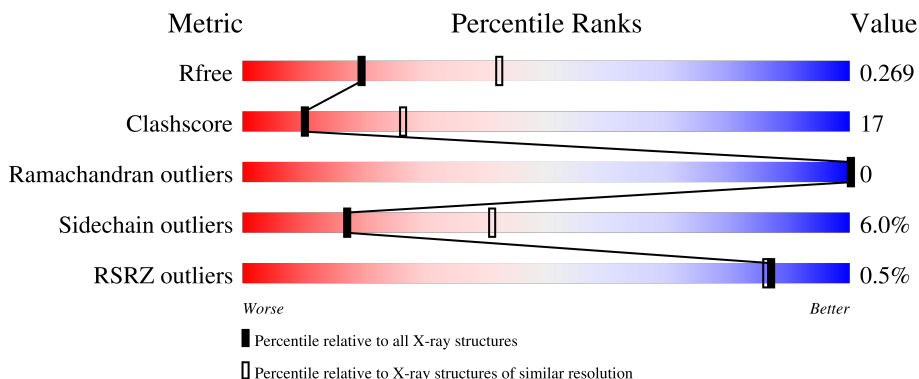
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	386	
1	B	386	
1	C	386	

2 Entry composition

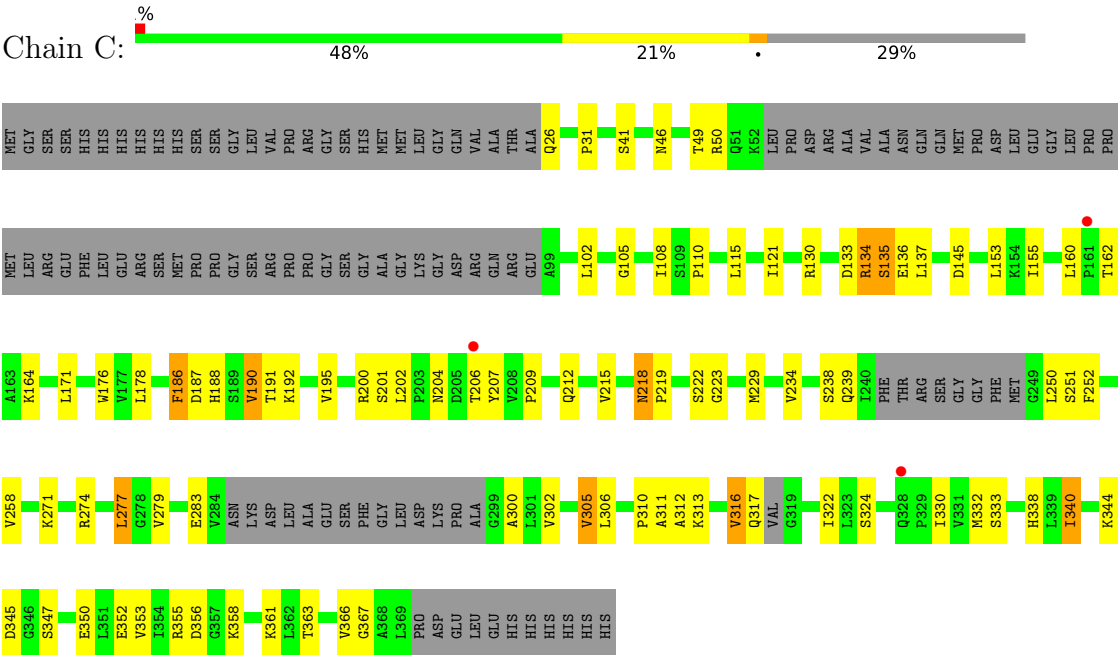
There is only 1 type of molecule in this entry. The entry contains 6193 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called serine endoprotease DegP-like protein MucD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	S	0	0	0
			2076	1302	360	411	3			
1	B	288	Total	C	N	O	S	0	0	0
			2108	1327	365	413	3			
1	C	275	Total	C	N	O	S	0	0	0
			2009	1263	350	393	3			

● Molecule 1: serine endoprotease DegP-like protein MucD



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	133.70Å 91.46Å 81.03Å 90.00° 116.06° 90.00°	Depositor
Resolution (Å)	29.56 – 2.70 29.56 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.2 (29.56-2.70) 99.3 (29.56-2.70)	Depositor EDS
R_{merge}	0.86	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 2.68Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.205 , 0.269 0.205 , 0.269	Depositor DCC
R_{free} test set	1998 reflections (8.23%)	wwPDB-VP
Wilson B-factor (Å ²)	62.8	Xtriage
Anisotropy	0.434	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 59.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.457 for -1/2*h+1/2*k+1,1/2*h-1/2*k+1,1/2*h+1/2*k 0.467 for -1/2*h-1/2*k+1,-1/2*h-1/2*k-1,1/2*h-1/2*k	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6193	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.97% 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

5.1 Standard geometry

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	0.48	0/2101	0.80	6/2852 (0.2%)
1	B	0.43	0/2135	0.74	1/2896 (0.0%)
1	C	0.41	0/2034	0.72	5/2759 (0.2%)
All	All	0.44	0/6270	0.75	12/8507 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
All	All	0	3

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	344	LYS	CG-CD-CE	-8.84	90.97	111.30
1	A	271	LYS	CA-CB-CG	8.83	131.75	114.10
1	A	274	ARG	CG-CD-NE	8.20	130.04	112.00
1	C	344	LYS	CB-CG-CD	7.38	128.27	111.30
1	B	200	ARG	CG-CD-NE	7.03	127.47	112.00
1	A	200	ARG	CD-NE-CZ	6.87	134.02	124.40
1	C	344	LYS	CD-CE-NZ	6.39	132.34	111.90
1	A	273	SER	CA-C-N	-6.02	112.02	122.37
1	A	273	SER	C-N-CA	-6.02	112.02	122.37
1	C	134	ARG	CB-CG-CD	5.09	123.02	111.30
1	A	313	LYS	CA-CB-CG	-5.04	104.02	114.10
1	C	134	ARG	CA-CB-CG	-5.00	104.10	114.10

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	200	ARG	Sidechain
1	A	274	ARG	Sidechain
1	B	200	ARG	Sidechain

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	2076	0	2108	72	0
1	B	2108	0	2152	77	0
1	C	2009	0	2052	69	0
All	All	6193	0	6312	214	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 17.

All (214) 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:324:SER:HB2	1:A:352:GLU:HB3	1.36	1.05
1:B:200:ARG:HA	1:B:209:PRO:HA	1.43	0.95
1:A:200:ARG:HA	1:A:209:PRO:HA	1.47	0.94
1:A:49:THR:HB	1:A:102:LEU:HD11	1.56	0.87
1:B:302:VAL:HG22	1:B:320:ASP:H	1.40	0.85
1:B:332:MET:HE1	1:B:334:ALA:HB3	1.60	0.84
1:C:305:VAL:HB	1:C:317:GLN:HG3	1.59	0.83
1:C:222:SER:HB2	1:C:239:GLN:HG2	1.67	0.76
1:C:200:ARG:HA	1:C:209:PRO:HA	1.68	0.76
1:C:218:ASN:HD22	1:C:219:PRO:HD2	1.51	0.74
1:B:209:PRO:HG2	1:B:258:VAL:HG21	1.72	0.71
1:A:28:GLU:HB2	1:C:229:MET:HB3	1.73	0.71
1:B:109:SER:OG	1:B:111:ASP:OD1	2.08	0.70
1:C:108:ILE:HD13	1:C:115:LEU:HG	1.73	0.70
1:C:134:ARG:NH1	1:C:186:PHE:HB2	2.08	0.68
1:A:48:SER:OG	1:A:130:ARG:NE	2.26	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:GLY:N	1:B:217:ILE:O	2.28	0.66
1:C:277:LEU:HD22	1:C:311:ALA:HB2	1.78	0.66
1:A:302:VAL:HG22	1:A:320:ASP:H	1.61	0.66
1:B:108:ILE:HD13	1:B:115:LEU:HG	1.77	0.66
1:B:280:VAL:HG13	1:B:304:GLN:HB2	1.78	0.65
1:C:209:PRO:HG2	1:C:258:VAL:HG21	1.79	0.64
1:C:352:GLU:HB3	1:C:361:LYS:HE3	1.80	0.64
1:B:200:ARG:HB3	1:B:209:PRO:HB3	1.80	0.64
1:C:41:SER:HB3	1:C:188:HIS:CD2	2.32	0.64
1:B:322:ILE:HD13	1:B:351:LEU:HD22	1.79	0.63
1:C:239:GLN:OE1	1:C:239:GLN:HA	1.97	0.63
1:B:279:VAL:HG21	1:B:302:VAL:HB	1.81	0.63
1:A:328:GLN:HB2	1:A:339:LEU:HD21	1.80	0.63
1:B:294:LEU:HD22	1:B:321:VAL:HG21	1.80	0.63
1:A:141:LEU:HD11	1:A:144:THR:HG22	1.81	0.62
1:B:27:ALA:O	1:B:28:GLU:HG2	2.00	0.62
1:C:218:ASN:O	1:C:239:GLN:NE2	2.25	0.61
1:B:322:ILE:HG12	1:B:353:VAL:HG12	1.81	0.61
1:A:50:ARG:HA	1:A:50:ARG:NH1	2.15	0.61
1:A:340:ILE:HA	1:A:343:LEU:HD22	1.83	0.61
1:C:49:THR:HB	1:C:102:LEU:HD11	1.82	0.61
1:C:145:ASP:OD2	1:C:274:ARG:NH2	2.33	0.61
1:C:207:TYR:OH	1:C:274:ARG:HD3	2.02	0.60
1:A:209:PRO:HG2	1:A:258:VAL:HG21	1.82	0.60
1:A:180:ILE:HG12	1:A:190:VAL:HG22	1.84	0.59
1:C:347:SER:O	1:C:366:VAL:HG12	2.02	0.59
1:A:238:SER:O	1:A:239:GLN:HG3	2.02	0.58
1:A:322:ILE:HG12	1:A:353:VAL:HG12	1.84	0.58
1:B:117:ASN:HD21	1:B:238:SER:HB2	1.69	0.58
1:A:28:GLU:HG3	1:A:30:LEU:H	1.68	0.58
1:C:356:ASP:HB2	1:C:358:LYS:HE2	1.86	0.57
1:A:37:VAL:O	1:A:41:SER:OG	2.19	0.57
1:A:108:ILE:HD13	1:A:115:LEU:HG	1.86	0.57
1:A:280:VAL:CG1	1:A:304:GLN:HB2	2.35	0.57
1:B:141:LEU:HD11	1:B:144:THR:CG2	2.35	0.56
1:C:324:SER:OG	1:C:352:GLU:HG2	2.06	0.56
1:A:348:LYS:HB3	1:A:363:THR:HG22	1.88	0.56
1:C:133:ASP:OD1	1:C:135:SER:OG	2.22	0.56
1:A:33:PHE:O	1:A:37:VAL:HG23	2.06	0.55
1:B:202:LEU:HD23	1:B:203:PRO:HG2	1.89	0.55
1:B:277:LEU:HD13	1:B:279:VAL:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:ILE:O	1:B:225:PRO:HD2	2.07	0.55
1:C:283:GLU:OE2	1:C:333:SER:N	2.29	0.55
1:A:352:GLU:HB2	1:A:361:LYS:HE2	1.89	0.54
1:B:352:GLU:HB2	1:B:361:LYS:HE3	1.88	0.54
1:C:277:LEU:HG	1:C:340:ILE:HD12	1.90	0.54
1:B:202:LEU:HD22	1:B:204:ASN:CG	2.33	0.54
1:B:31:PRO:HB3	1:C:26:GLN:HB3	1.90	0.54
1:A:265:GLN:OE1	1:A:272:VAL:HG13	2.09	0.53
1:A:322:ILE:HG22	1:A:330:ILE:HD12	1.90	0.53
1:A:317:GLN:N	1:A:317:GLN:OE1	2.38	0.53
1:B:141:LEU:HD11	1:B:144:THR:HG22	1.90	0.53
1:B:355:ARG:HH22	1:B:362:LEU:HD21	1.73	0.53
1:B:33:PHE:O	1:B:37:VAL:HG23	2.08	0.53
1:A:46:ASN:OD1	1:A:130:ARG:NE	2.42	0.53
1:B:212:GLN:OE1	1:B:250:LEU:HD13	2.09	0.53
1:B:49:THR:HB	1:B:102:LEU:HD11	1.91	0.52
1:B:283:GLU:OE2	1:B:332:MET:HG2	2.09	0.52
1:A:180:ILE:O	1:A:225:PRO:HD2	2.10	0.52
1:C:46:ASN:OD1	1:C:130:ARG:HG3	2.08	0.52
1:B:199:GLY:C	1:B:200:ARG:HG2	2.31	0.52
1:B:337:PRO:O	1:B:341:GLY:N	2.39	0.52
1:C:317:GLN:H	1:C:317:GLN:CD	2.16	0.52
1:A:239:GLN:HA	1:A:252:PHE:H	1.76	0.51
1:C:218:ASN:H	1:C:239:GLN:NE2	2.08	0.51
1:B:302:VAL:HG23	1:B:318:VAL:HA	1.93	0.51
1:B:277:LEU:HD12	1:B:278:GLY:H	1.75	0.51
1:B:355:ARG:O	1:B:358:LYS:HG2	2.11	0.51
1:B:182:SER:HB2	1:B:188:HIS:HD2	1.76	0.51
1:A:202:LEU:HD12	1:A:204:ASN:OD1	2.11	0.51
1:A:324:SER:OG	1:A:352:GLU:OE1	2.12	0.51
1:A:202:LEU:HD13	1:A:203:PRO:HD2	1.93	0.51
1:B:49:THR:HA	1:B:126:GLU:O	2.11	0.51
1:C:355:ARG:HB3	1:C:358:LYS:HE3	1.92	0.50
1:C:110:PRO:N	1:C:162:THR:HG21	2.26	0.50
1:A:324:SER:CB	1:A:352:GLU:HB3	2.24	0.50
1:A:273:SER:C	1:A:274:ARG:HG2	2.37	0.50
1:B:305:VAL:HG11	1:B:316:VAL:O	2.12	0.50
1:A:273:SER:O	1:A:274:ARG:HG2	2.12	0.50
1:C:279:VAL:HG12	1:C:305:VAL:HG22	1.94	0.49
1:A:348:LYS:HA	1:A:364:VAL:O	2.11	0.49
1:B:48:SER:OG	1:B:130:ARG:HD2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:ARG:HA	1:A:50:ARG:HH11	1.76	0.49
1:C:306:LEU:O	1:C:312:ALA:HB2	2.13	0.49
1:C:322:ILE:HG12	1:C:353:VAL:HG12	1.95	0.49
1:A:330:ILE:HG23	1:A:332:MET:O	2.12	0.49
1:B:328:GLN:HG3	1:B:339:LEU:HD11	1.95	0.49
1:C:171:LEU:HG	1:C:195:VAL:HG21	1.94	0.49
1:B:317:GLN:NE2	1:B:320:ASP:OD1	2.45	0.48
1:B:134:ARG:HA	1:B:186:PHE:CD2	2.47	0.48
1:C:176:TRP:C	1:C:229:MET:HE3	2.38	0.48
1:B:205:ASP:OD1	1:B:206:THR:N	2.46	0.48
1:A:115:LEU:HD13	1:A:236:ILE:HD11	1.96	0.48
1:C:316:VAL:HG13	1:C:317:GLN:N	2.28	0.48
1:B:277:LEU:HD12	1:B:278:GLY:N	2.28	0.48
1:A:339:LEU:O	1:A:343:LEU:HD13	2.14	0.48
1:B:332:MET:HE3	1:B:335:ASP:OD1	2.13	0.48
1:A:345:ASP:OD1	1:A:367:GLY:HA2	2.14	0.48
1:B:176:TRP:HB3	1:B:229:MET:HE3	1.96	0.48
1:C:121:ILE:HG21	1:C:153:LEU:HD21	1.94	0.48
1:C:300:ALA:HB2	1:C:333:SER:HA	1.96	0.48
1:A:310:PRO:HB2	1:A:364:VAL:HG13	1.96	0.47
1:C:178:LEU:HD11	1:C:190:VAL:HG13	1.96	0.47
1:C:271:LYS:HB3	1:C:271:LYS:HE2	1.72	0.47
1:A:40:ALA:HB1	1:A:227:PHE:CZ	2.49	0.47
1:A:115:LEU:HD11	1:A:260:MET:CE	2.44	0.47
1:A:192:LYS:O	1:A:192:LYS:HD3	2.15	0.47
1:A:202:LEU:HD22	1:A:203:PRO:HD2	1.97	0.47
1:C:50:ARG:HD2	1:C:50:ARG:HA	1.69	0.47
1:A:196:SER:N	1:A:212:GLN:O	2.43	0.47
1:C:238:SER:O	1:C:252:PHE:O	2.33	0.47
1:B:289:ALA:HA	1:B:294:LEU:HB3	1.96	0.47
1:B:212:GLN:HG3	1:B:252:PHE:CE1	2.49	0.46
1:B:349:ALA:HB3	1:B:364:VAL:HB	1.97	0.46
1:C:137:LEU:HD12	1:C:155:ILE:HD12	1.97	0.46
1:A:106:PHE:HB3	1:A:225:PRO:HG3	1.97	0.46
1:C:191:THR:HG23	1:C:215:VAL:HG13	1.96	0.46
1:A:290:GLU:OE1	1:A:291:SER:N	2.49	0.46
1:A:202:LEU:HD13	1:A:203:PRO:CD	2.45	0.46
1:C:136:GLU:O	1:C:137:LEU:HD23	2.15	0.46
1:A:192:LYS:HD3	1:A:192:LYS:C	2.41	0.45
1:B:150:VAL:HB	1:B:262:VAL:HG11	1.99	0.45
1:B:307:GLU:C	1:B:309:GLY:H	2.23	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:GLU:CD	1:B:332:MET:HG2	2.42	0.45
1:B:332:MET:HE3	1:B:332:MET:HB3	1.62	0.45
1:C:316:VAL:HG12	1:C:355:ARG:NH2	2.31	0.45
1:C:332:MET:HE2	1:C:332:MET:HB2	1.92	0.45
1:B:280:VAL:CG1	1:B:304:GLN:HB2	2.46	0.45
1:B:171:LEU:HG	1:B:195:VAL:HG21	1.98	0.45
1:C:316:VAL:HG22	1:C:317:GLN:OE1	2.17	0.45
1:B:354:ILE:HD13	1:B:359:ARG:HG3	1.99	0.45
1:B:116:THR:CG2	1:B:151:ALA:HB3	2.47	0.44
1:B:209:PRO:CG	1:B:258:VAL:HG21	2.45	0.44
1:C:345:ASP:OD1	1:C:367:GLY:HA2	2.18	0.44
1:B:143:GLY:C	1:B:266:LEU:HD13	2.41	0.44
1:B:344:LYS:HD2	1:B:344:LYS:HA	1.68	0.44
1:A:336:LEU:HB3	1:A:337:PRO:HD3	1.98	0.44
1:B:46:ASN:HB2	1:B:182:SER:OG	2.18	0.44
1:B:119:HIS:ND1	1:B:149:ASP:OD2	2.43	0.44
1:C:105:GLY:HA2	1:C:223:GLY:O	2.17	0.44
1:B:300:ALA:HB2	1:B:333:SER:HA	2.00	0.44
1:C:350:GLU:HG3	1:C:363:THR:OG1	2.18	0.44
1:B:354:ILE:CD1	1:B:359:ARG:HG3	2.48	0.44
1:C:283:GLU:OE2	1:C:332:MET:HG3	2.18	0.43
1:A:332:MET:O	1:A:335:ASP:HB2	2.18	0.43
1:B:302:VAL:HG13	1:B:320:ASP:O	2.18	0.43
1:C:171:LEU:HD22	1:C:234:VAL:HB	1.99	0.43
1:A:36:LEU:HD23	1:A:178:LEU:HD22	1.99	0.43
1:B:279:VAL:HB	1:B:305:VAL:HA	2.01	0.43
1:B:334:ALA:O	1:B:338:HIS:HB2	2.19	0.43
1:B:355:ARG:HD2	1:B:358:LYS:HE3	2.01	0.43
1:A:141:LEU:HD11	1:A:144:THR:CG2	2.49	0.43
1:B:40:ALA:HB1	1:B:227:PHE:CZ	2.54	0.43
1:C:352:GLU:O	1:C:352:GLU:HG3	2.19	0.43
1:A:171:LEU:HG	1:A:195:VAL:HG21	2.00	0.42
1:B:176:TRP:C	1:B:229:MET:HE3	2.44	0.42
1:B:176:TRP:NE1	1:C:31:PRO:O	2.40	0.42
1:B:183:PRO:HB3	1:B:220:GLY:HA3	2.01	0.42
1:C:160:LEU:O	1:C:162:THR:HG23	2.19	0.42
1:C:134:ARG:CZ	1:C:186:PHE:HB2	2.49	0.42
1:A:333:SER:O	1:A:337:PRO:HD3	2.19	0.42
1:B:126:GLU:OE2	1:B:127:ILE:N	2.53	0.42
1:C:164:LYS:HB2	1:C:164:LYS:HE3	1.71	0.42
1:C:212:GLN:OE1	1:C:250:LEU:HD13	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:ALA:O	1:A:337:PRO:HD2	2.20	0.42
1:A:266:LEU:CD1	1:A:272:VAL:HG23	2.50	0.42
1:B:332:MET:CE	1:B:334:ALA:HB3	2.42	0.42
1:C:302:VAL:HG11	1:C:305:VAL:HG23	2.02	0.41
1:A:278:GLY:O	1:A:306:LEU:HG	2.19	0.41
1:A:317:GLN:HE22	1:A:355:ARG:HB2	1.86	0.41
1:A:348:LYS:HE3	1:A:365:THR:OG1	2.20	0.41
1:A:276:TRP:HD1	1:A:367:GLY:C	2.28	0.41
1:C:134:ARG:NH2	1:C:187:ASP:OD1	2.53	0.41
1:A:294:LEU:HD12	1:A:294:LEU:HA	1.88	0.41
1:C:212:GLN:HG3	1:C:252:PHE:CE2	2.56	0.41
1:A:28:GLU:HG3	1:A:30:LEU:HB2	2.02	0.41
1:A:321:VAL:HB	1:A:354:ILE:HG22	2.02	0.41
1:A:104:SER:OG	1:A:220:GLY:O	2.28	0.41
1:C:204:ASN:OD1	1:C:204:ASN:N	2.48	0.41
1:A:130:ARG:HB3	1:A:186:PHE:CD2	2.56	0.41
1:A:130:ARG:HB3	1:A:186:PHE:CE2	2.55	0.41
1:C:277:LEU:HD23	1:C:277:LEU:HA	1.78	0.41
1:C:302:VAL:HG11	1:C:317:GLN:HA	2.03	0.41
1:A:41:SER:HB3	1:A:188:HIS:ND1	2.36	0.41
1:A:302:VAL:HG13	1:A:320:ASP:O	2.20	0.41
1:B:136:GLU:O	1:B:137:LEU:HD23	2.21	0.41
1:C:145:ASP:CG	1:C:274:ARG:HH21	2.29	0.40
1:C:322:ILE:HG22	1:C:330:ILE:HD12	2.03	0.40
1:B:142:VAL:HB	1:B:152:VAL:HG12	2.04	0.40
1:B:279:VAL:HG23	1:B:304:GLN:O	2.22	0.40
1:C:202:LEU:HB3	1:C:204:ASN:OD1	2.21	0.40
1:A:176:TRP:C	1:A:229:MET:HE3	2.46	0.40
1:C:191:THR:CG2	1:C:215:VAL:HG13	2.51	0.40
1:C:330:ILE:HG23	1:C:332:MET:O	2.21	0.40
1:A:347:SER:O	1:A:366:VAL:HG22	2.22	0.40
1:B:137:LEU:HD12	1:B:155:ILE:HD12	2.04	0.40
1:B:217:ILE:HD12	1:B:217:ILE:H	1.87	0.40
1:C:310:PRO:O	1:C:313:LYS:HB3	2.22	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	278/386 (72%)	265 (95%)	13 (5%)	0	100	100
1	B	280/386 (72%)	263 (94%)	17 (6%)	0	100	100
1	C	267/386 (69%)	255 (96%)	12 (4%)	0	100	100
All	All	825/1158 (71%)	783 (95%)	42 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	225/310 (73%)	212 (94%)	13 (6%)	18	42
1	B	229/310 (74%)	215 (94%)	14 (6%)	17	40
1	C	218/310 (70%)	205 (94%)	13 (6%)	17	41
All	All	672/930 (72%)	632 (94%)	40 (6%)	17	41

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	SER
1	A	48	SER
1	A	187	ASP
1	A	190	VAL
1	A	192	LYS

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Mol	Chain	Res	Type
1	A	200	ARG
1	A	201	SER
1	A	202	LEU
1	A	215	VAL
1	A	240	ILE
1	A	307	GLU
1	A	318	VAL
1	A	362	LEU
1	B	41	SER
1	B	48	SER
1	B	126	GLU
1	B	190	VAL
1	B	192	LYS
1	B	200	ARG
1	B	256	ILE
1	B	277	LEU
1	B	279	VAL
1	B	306	LEU
1	B	313	LYS
1	B	318	VAL
1	B	338	HIS
1	B	343	LEU
1	C	135	SER
1	C	186	PHE
1	C	190	VAL
1	C	192	LYS
1	C	201	SER
1	C	206	THR
1	C	218	ASN
1	C	251	SER
1	C	277	LEU
1	C	305	VAL
1	C	316	VAL
1	C	338	HIS
1	C	340	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	328	GLN
1	B	169	ASN
1	B	282	GLN

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Mol	Chain	Res	Type
1	C	29	ASN
1	C	39	GLN
1	C	169	ASN
1	C	218	ASN
1	C	221	ASN
1	C	237	ASN
1	C	265	GLN
1	C	282	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 [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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	286/386 (74%)	-0.89	0 100 100	39, 71, 131, 144	0
1	B	288/386 (74%)	-0.87	1 (0%) 90 89	39, 72, 128, 146	0
1	C	275/386 (71%)	-0.87	3 (1%) 78 76	41, 68, 123, 151	0
All	All	849/1158 (73%)	-0.88	4 (0%) 87 86	39, 70, 129, 151	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	328	GLN	3.8
1	C	206	THR	2.2
1	C	161	PRO	2.1
1	B	252	PHE	2.1

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.