



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:58 AM GMT

PDB ID : 3CUQ
Title : Integrated structural and functional model of the human ESCRT-II complex
Authors : Im, Y.J.; Hurley, J.H.
Deposited on : 2008-04-16
Resolution : 2.61 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

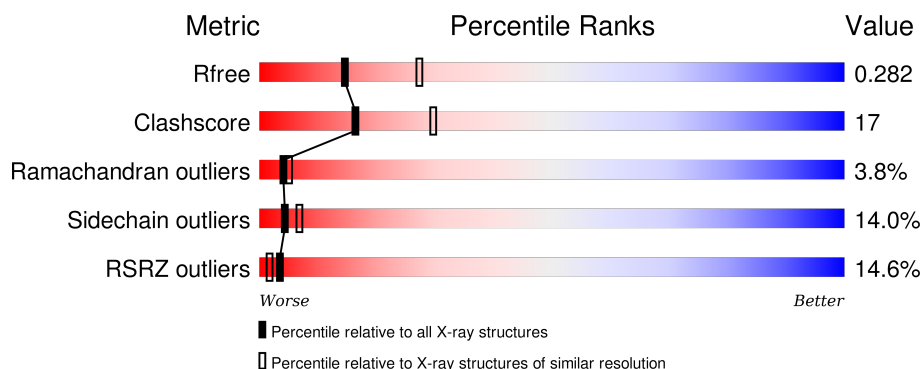
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 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	2700 (2.64-2.60)
Clashscore	102246	3065 (2.64-2.60)
Ramachandran outliers	100387	3015 (2.64-2.60)
Sidechain outliers	100360	3015 (2.64-2.60)
RSRZ outliers	91569	2706 (2.64-2.60)

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	234	<div> <div>12%</div> <div>62%</div> <div>25%</div> <div>6%</div> <div>6%</div> </div>
2	B	218	<div> <div>14%</div> <div>62%</div> <div>27%</div> <div>5%</div> <div>6%</div> </div>
3	C	176	<div> <div>6%</div> <div>49%</div> <div>40%</div> <div>9%</div> <div>• •</div> </div>
3	D	176	<div> <div>19%</div> <div>31%</div> <div>18%</div> <div>5%</div> <div>•</div> <div>45%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5630 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 Vacuolar-sorting protein SNF8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1736	1113	292	324	7			

- Molecule 2 is a protein called Vacuolar protein-sorting-associated protein 36.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	204	Total	C	N	O	S	0	0	0
			1593	1007	267	304	15			

- Molecule 3 is a protein called Vacuolar protein-sorting-associated protein 25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	173	Total	C	N	O	S	0	0	0
			1442	926	247	264	5			
3	D	97	Total	C	N	O	S	0	0	0
			821	537	139	140	5			

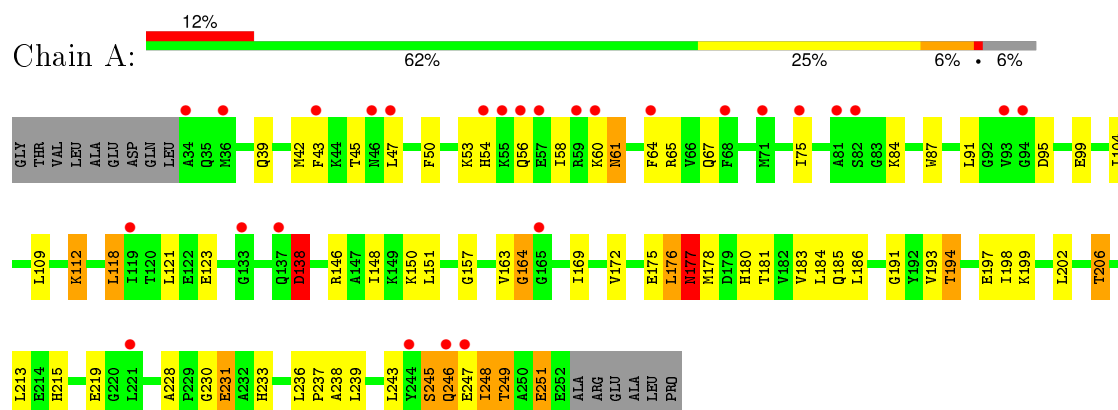
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	O	0	0
			12	12		
4	B	13	Total	O	0	0
			13	13		
4	C	10	Total	O	0	0
			10	10		
4	D	3	Total	O	0	0
			3	3		

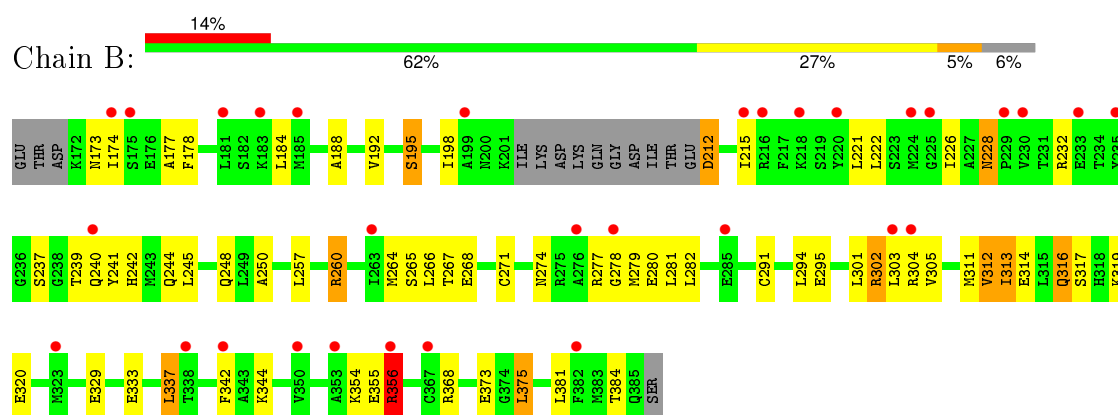
3 Residue-property plots

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.

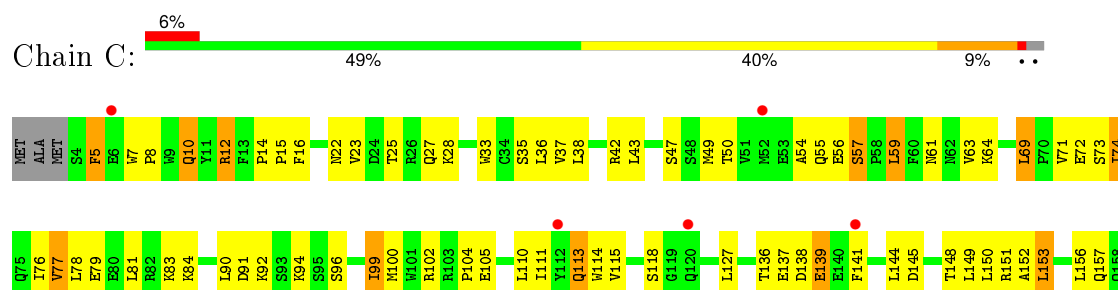
• Molecule 1: Vacuolar-sorting protein SNF8

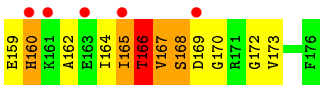


• Molecule 2: Vacuolar protein-sorting-associated protein 36

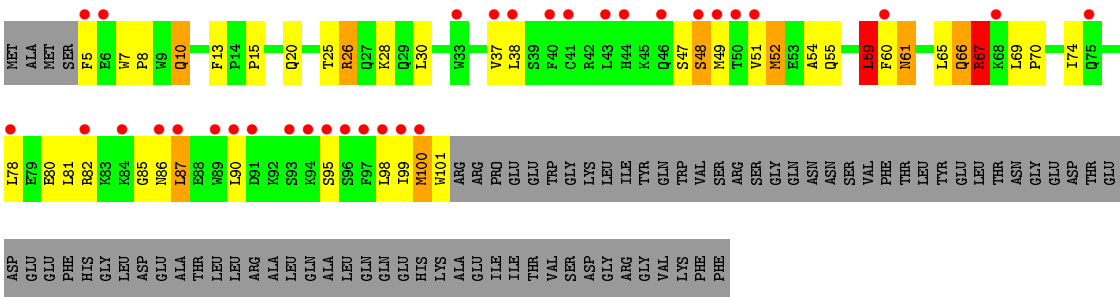
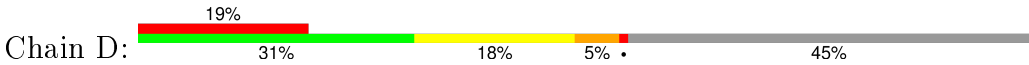


• Molecule 3: Vacuolar protein-sorting-associated protein 25





● Molecule 3: Vacuolar protein-sorting-associated protein 25



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.16Å 89.15Å 91.44Å 90.00° 101.52° 90.00°	Depositor
Resolution (Å)	44.80 – 2.61 44.80 – 2.61	Depositor EDS
% Data completeness (in resolution range)	91.0 (44.80-2.61) 91.0 (44.80-2.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.10 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.238 , 0.298 0.226 , 0.282	Depositor DCC
R_{free} test set	1532 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	69.0	Xtriage
Anisotropy	0.638	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 93.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 30665 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5630	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% 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	0.76	2/1770 (0.1%)	0.83	0/2389
2	B	0.72	2/1610 (0.1%)	0.87	4/2157 (0.2%)
3	C	0.70	0/1479	0.78	0/2000
3	D	0.51	0/845	0.66	1/1143 (0.1%)
All	All	0.70	4/5704 (0.1%)	0.81	5/7689 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	177	ASN	CB-CG	6.12	1.65	1.51
1	A	123	GLU	CG-CD	6.07	1.61	1.51
2	B	271	CYS	CB-SG	-5.57	1.72	1.81
2	B	355	GLU	CG-CD	5.08	1.59	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	337	LEU	CB-CG-CD1	-8.22	97.03	111.00
2	B	356	ARG	NE-CZ-NH1	-7.01	116.79	120.30
2	B	368	ARG	NE-CZ-NH1	-5.61	117.49	120.30
3	D	59	LEU	CA-CB-CG	5.23	127.33	115.30
2	B	294	LEU	CA-CB-CG	5.15	127.14	115.30

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	1736	0	1742	45	0
2	B	1593	0	1649	60	0
3	C	1442	0	1416	66	0
3	D	821	0	828	24	0
4	A	12	0	0	1	0
4	B	13	0	0	1	0
4	C	10	0	0	0	0
4	D	3	0	0	0	0
All	All	5630	0	5635	187	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 (187) 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:B:302:ARG:HH21	2:B:304:ARG:HG2	1.18	1.06
3:C:100:MET:HE1	3:C:104:PRO:HG3	1.40	1.00
3:C:167:VAL:HG22	3:C:168:SER:H	1.30	0.96
3:C:100:MET:CE	3:C:104:PRO:HG3	1.97	0.94
3:C:10:GLN:HE21	3:C:10:GLN:H	1.15	0.93
2:B:228:ASN:H	2:B:232:ARG:HD3	1.40	0.86
3:C:37:VAL:HG11	3:C:81:LEU:HD11	1.59	0.85
1:A:194:THR:HG22	1:A:197:GLU:H	1.43	0.83
3:C:136:THR:O	3:C:138:ASP:N	2.12	0.82
3:C:165:ILE:HG22	3:C:166:THR:N	1.95	0.82
1:A:199:LYS:NZ	1:A:206:THR:HB	1.96	0.79
2:B:245:LEU:HD21	2:B:282:LEU:HD11	1.64	0.78
2:B:295:GLU:OE2	2:B:302:ARG:HG3	1.82	0.78
3:D:20:GLN:O	3:D:26:ARG:HD3	1.84	0.77
3:C:153:LEU:HD23	3:C:173:VAL:HG21	1.66	0.77
2:B:212:ASP:HA	2:B:215:ILE:HD12	1.68	0.75
3:C:14:PRO:HB2	3:C:15:PRO:HD3	1.70	0.74
1:A:104:ILE:HG21	1:A:148:ILE:HD11	1.69	0.73
2:B:337:LEU:HD12	2:B:337:LEU:C	2.07	0.73
1:A:138:ASP:N	1:A:138:ASP:OD1	2.21	0.72
3:C:165:ILE:HG22	3:C:166:THR:H	1.53	0.72
2:B:337:LEU:CD1	2:B:342:PHE:HB2	2.21	0.71
2:B:302:ARG:HH21	2:B:304:ARG:CG	1.99	0.70
2:B:320:GLU:OE1	2:B:356:ARG:HD3	1.92	0.69
3:C:111:ILE:O	3:C:115:VAL:HG23	1.94	0.68
2:B:274:ASN:HD21	2:B:282:LEU:H	1.42	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:37:VAL:CG1	3:C:81:LEU:HD11	2.25	0.67
2:B:274:ASN:ND2	2:B:282:LEU:H	1.92	0.66
3:C:172:GLY:O	3:C:173:VAL:HG23	1.96	0.66
3:D:66:GLN:O	3:D:67:ARG:HG2	1.96	0.66
2:B:291:CYS:SG	2:B:313:ILE:HD11	2.36	0.65
1:A:199:LYS:HZ2	1:A:206:THR:HB	1.61	0.64
2:B:337:LEU:HD12	2:B:337:LEU:O	1.97	0.64
2:B:337:LEU:HD11	2:B:342:PHE:HB2	1.78	0.64
3:D:52:MET:HA	3:D:55:GLN:CD	2.16	0.64
2:B:373:GLU:O	3:C:23:VAL:HG12	1.98	0.64
3:C:100:MET:HE3	3:C:148:THR:HG23	1.78	0.63
3:D:87:LEU:HA	3:D:98:LEU:O	1.99	0.63
1:A:163:VAL:O	1:A:164:GLY:C	2.37	0.63
2:B:239:THR:HG22	2:B:241:TYR:H	1.64	0.62
3:D:81:LEU:O	3:D:85:GLY:O	2.18	0.61
3:C:10:GLN:NE2	3:C:10:GLN:H	1.93	0.60
1:A:104:ILE:HG21	1:A:148:ILE:CD1	2.31	0.60
2:B:245:LEU:O	2:B:248:GLN:HB2	2.02	0.59
3:C:22:ASN:ND2	3:C:25:THR:H	2.01	0.59
3:C:167:VAL:CG2	3:C:168:SER:H	2.12	0.58
2:B:265:SER:HA	2:B:312:VAL:HG23	1.86	0.58
3:D:20:GLN:HG3	3:D:25:THR:CG2	2.34	0.58
2:B:381:LEU:HD12	3:D:13:PHE:CE1	2.38	0.58
2:B:228:ASN:N	2:B:232:ARG:HD3	2.16	0.57
3:C:69:LEU:HD22	3:C:73:SER:HB2	1.86	0.57
3:C:167:VAL:HG22	3:C:168:SER:N	2.12	0.57
1:A:177:ASN:OD1	1:A:180:HIS:HB2	2.05	0.57
1:A:199:LYS:HZ3	1:A:206:THR:HB	1.67	0.56
2:B:257:LEU:HD23	2:B:264:MET:HG3	1.87	0.56
1:A:237:PRO:O	1:A:239:LEU:O	2.24	0.56
3:D:20:GLN:HG3	3:D:25:THR:HG22	1.88	0.55
2:B:266:LEU:HD23	2:B:313:ILE:HG13	1.88	0.55
2:B:274:ASN:HD21	2:B:282:LEU:N	2.04	0.55
2:B:239:THR:HB	2:B:242:HIS:CD2	2.41	0.55
2:B:195:SER:HA	2:B:198:ILE:HG12	1.89	0.55
2:B:228:ASN:H	2:B:232:ARG:HH11	1.54	0.55
3:D:47:SER:O	3:D:48:SER:HB2	2.07	0.55
2:B:337:LEU:HD13	2:B:342:PHE:HB2	1.88	0.54
3:C:100:MET:HE2	3:C:104:PRO:HG3	1.84	0.54
3:C:166:THR:O	3:C:170:GLY:O	2.25	0.54
1:A:181:THR:O	1:A:185:GLN:HG3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:5:PHE:C	3:C:5:PHE:CD2	2.81	0.53
2:B:274:ASN:HB3	2:B:281:LEU:HD13	1.90	0.53
3:C:71:VAL:O	3:C:74:ILE:HG12	2.08	0.53
2:B:274:ASN:HD21	2:B:282:LEU:HB2	1.74	0.52
1:A:194:THR:HG22	1:A:197:GLU:N	2.17	0.52
2:B:316:GLN:HE21	2:B:316:GLN:HA	1.73	0.52
3:C:100:MET:CE	3:C:148:THR:HG23	2.39	0.52
3:D:51:VAL:HG12	3:D:55:GLN:HE21	1.76	0.51
3:C:74:ILE:HA	3:C:77:VAL:CG1	2.40	0.51
3:C:153:LEU:CD2	3:C:173:VAL:HG21	2.39	0.51
2:B:195:SER:HA	2:B:198:ILE:CG1	2.41	0.51
3:D:37:VAL:HG11	3:D:81:LEU:CD2	2.41	0.51
3:C:37:VAL:HG11	3:C:81:LEU:CD1	2.36	0.50
1:A:163:VAL:O	1:A:164:GLY:O	2.28	0.50
3:D:37:VAL:HG11	3:D:81:LEU:HD21	1.92	0.50
3:C:150:LEU:O	3:C:151:ARG:C	2.49	0.50
3:D:61:ASN:HD22	3:D:61:ASN:C	2.15	0.50
1:A:87:TRP:O	1:A:91:LEU:HB2	2.10	0.50
2:B:375:LEU:C	2:B:375:LEU:HD12	2.32	0.50
2:B:260:ARG:HH11	2:B:260:ARG:HG3	1.76	0.49
2:B:264:MET:CE	2:B:268:GLU:HB3	2.41	0.49
3:C:69:LEU:HD13	3:C:74:ILE:CG2	2.42	0.49
3:C:61:ASN:OD1	3:C:63:VAL:HG22	2.12	0.49
3:C:47:SER:HB3	3:C:145:ASP:OD1	2.12	0.49
2:B:260:ARG:NH1	2:B:260:ARG:HG3	2.27	0.49
2:B:250:ALA:HA	2:B:301:LEU:HD11	1.94	0.49
3:C:57:SER:HB2	3:C:59:LEU:H	1.76	0.49
1:A:194:THR:O	1:A:198:ILE:HG12	2.12	0.49
3:D:82:ARG:HA	3:D:87:LEU:HB3	1.95	0.49
2:B:244:GLN:OE1	2:B:244:GLN:HA	2.13	0.49
1:A:95:ASP:O	1:A:99:GLU:HB2	2.12	0.49
2:B:305:VAL:HG22	2:B:311:MET:HE2	1.95	0.49
1:A:176:LEU:O	1:A:177:ASN:C	2.51	0.48
3:C:102:ARG:NH1	3:C:110:LEU:HD11	2.28	0.48
1:A:84:LYS:HD3	2:B:222:LEU:HD23	1.93	0.48
3:C:113:GLN:HE21	3:C:114:TRP:N	2.10	0.48
3:C:5:PHE:O	3:C:5:PHE:CD2	2.67	0.48
2:B:302:ARG:HE	2:B:304:ARG:HG3	1.78	0.48
3:D:52:MET:HA	3:D:55:GLN:NE2	2.28	0.48
3:C:136:THR:HG23	3:C:139:GLU:HB2	1.96	0.48
3:D:30:LEU:HD22	3:D:80:GLU:HG3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:13:PHE:CD2	3:D:15:PRO:HD2	2.49	0.47
3:C:74:ILE:HA	3:C:77:VAL:HG13	1.97	0.47
2:B:277:ARG:HH12	2:B:282:LEU:HD13	1.79	0.47
3:C:22:ASN:HD22	3:C:25:THR:H	1.61	0.47
1:A:43:PHE:CZ	2:B:188:ALA:HB2	2.50	0.47
1:A:230:GLY:O	1:A:231:GLU:HB3	2.14	0.47
1:A:109:LEU:HD23	1:A:112:LYS:HE3	1.96	0.47
1:A:177:ASN:H	1:A:177:ASN:HD22	1.64	0.46
3:C:8:PRO:HD3	3:C:36:LEU:HD13	1.97	0.46
2:B:221:LEU:HD22	2:B:226:ILE:HB	1.95	0.46
3:C:172:GLY:O	3:C:173:VAL:CG2	2.64	0.46
2:B:257:LEU:CD2	2:B:264:MET:HG3	2.45	0.45
3:C:157:GLN:HA	3:C:162:ALA:O	2.16	0.45
1:A:75:ILE:HG22	2:B:192:VAL:HG21	1.97	0.45
3:C:73:SER:O	3:C:76:ILE:HB	2.16	0.45
2:B:304:ARG:HH22	2:B:320:GLU:CG	2.30	0.45
3:C:165:ILE:CG2	3:C:166:THR:H	2.19	0.45
3:C:5:PHE:O	3:C:5:PHE:HD2	2.00	0.45
2:B:320:GLU:OE1	2:B:356:ARG:CD	2.62	0.45
3:D:7:TRP:HA	3:D:8:PRO:HD3	1.86	0.45
3:C:114:TRP:O	3:C:118:SER:HB2	2.17	0.45
3:C:10:GLN:N	3:C:10:GLN:HE21	1.98	0.45
3:C:71:VAL:HA	3:C:74:ILE:HD11	1.98	0.45
1:A:177:ASN:N	1:A:177:ASN:HD22	2.14	0.45
3:C:72:GLU:HG2	3:C:73:SER:N	2.33	0.44
3:C:54:ALA:O	3:C:56:GLU:N	2.50	0.44
1:A:247:GLU:O	1:A:248:ILE:HB	2.17	0.44
3:C:69:LEU:HD13	3:C:74:ILE:HG22	2.00	0.44
3:C:56:GLU:HA	3:C:56:GLU:OE1	2.17	0.44
3:C:37:VAL:HG13	3:C:38:LEU:H	1.83	0.44
3:D:86:ASN:HA	3:D:100:MET:HB2	1.99	0.43
1:A:67:GLN:HE21	2:B:178:PHE:HE2	1.65	0.43
1:A:39:GLN:HA	1:A:42:MET:HG3	2.00	0.43
1:A:47:LEU:O	1:A:50:PHE:HB3	2.18	0.43
3:D:54:ALA:HB1	3:D:60:PHE:HD1	1.83	0.43
3:C:166:THR:O	3:C:167:VAL:O	2.37	0.43
3:C:111:ILE:HG12	3:C:141:PHE:CD2	2.53	0.43
2:B:304:ARG:HD2	4:B:23:HOH:O	2.19	0.43
3:C:57:SER:C	3:C:59:LEU:H	2.22	0.43
1:A:246:GLN:O	1:A:246:GLN:HG2	2.18	0.43
1:A:180:HIS:HE1	1:A:184:LEU:HD11	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:LEU:HA	1:A:237:PRO:HD2	1.85	0.43
3:C:16:PHE:CE2	3:C:33:TRP:CE2	3.07	0.42
3:C:43:LEU:C	3:C:43:LEU:HD13	2.40	0.42
2:B:279:MET:O	2:B:280:GLU:HG2	2.19	0.42
2:B:266:LEU:CD2	2:B:313:ILE:HG13	2.48	0.42
3:C:47:SER:CB	3:C:145:ASP:OD1	2.67	0.42
1:A:118:LEU:HB3	1:A:238:ALA:HB2	2.01	0.42
3:D:59:LEU:HD13	3:D:60:PHE:N	2.34	0.42
1:A:215:HIS:NE2	1:A:219:GLU:OE1	2.52	0.42
3:C:79:GLU:O	3:C:83:LYS:HG2	2.19	0.42
3:C:71:VAL:HA	3:C:74:ILE:CD1	2.49	0.42
1:A:56:GLN:O	1:A:60:LYS:HB2	2.20	0.42
1:A:58:ILE:HG23	1:A:65:ARG:HG2	2.01	0.42
3:C:99:ILE:O	3:C:99:ILE:HG23	2.20	0.42
1:A:61:ASN:HD21	1:A:64:PHE:CB	2.33	0.42
1:A:249:THR:HA	1:A:251:GLU:OE2	2.20	0.41
3:C:7:TRP:HB2	3:C:12:ARG:HH12	1.85	0.41
1:A:191:GLY:HA2	1:A:239:LEU:CD1	2.50	0.41
3:C:14:PRO:HB2	3:C:15:PRO:CD	2.46	0.41
2:B:239:THR:HG22	2:B:240:GLN:N	2.35	0.41
1:A:53:LYS:HE2	2:B:173:ASN:HD21	1.84	0.41
2:B:291:CYS:HB3	2:B:303:LEU:HD22	2.02	0.41
1:A:54:HIS:CD2	2:B:177:ALA:HB2	2.55	0.41
1:A:151:LEU:HA	1:A:151:LEU:HD23	1.84	0.41
3:C:149:LEU:O	3:C:152:ALA:HB3	2.21	0.41
2:B:264:MET:O	2:B:312:VAL:HG22	2.21	0.41
1:A:228:ALA:HB1	1:A:233:HIS:CE1	2.56	0.41
3:D:100:MET:HB3	3:D:101:TRP:H	1.55	0.41
3:D:70:PRO:O	3:D:74:ILE:HG12	2.20	0.41
2:B:274:ASN:ND2	2:B:282:LEU:N	2.66	0.41
1:A:183:VAL:O	1:A:184:LEU:C	2.59	0.41
1:A:243:LEU:HD21	4:A:263:HOH:O	2.22	0.41
2:B:329:GLU:OE1	2:B:333:GLU:OE2	2.40	0.41
2:B:302:ARG:NH2	2:B:304:ARG:HG2	2.04	0.40
1:A:121:LEU:HD13	1:A:169:ILE:HD11	2.04	0.40
3:C:23:VAL:O	3:C:27:GLN:HG3	2.22	0.40
3:D:10:GLN:HE21	3:D:10:GLN:H	1.69	0.40
2:B:337:LEU:CD1	2:B:337:LEU:C	2.82	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	217/234 (93%)	197 (91%)	13 (6%)	7 (3%)	5	7
2	B	200/218 (92%)	182 (91%)	15 (8%)	3 (2%)	13	24
3	C	171/176 (97%)	145 (85%)	16 (9%)	10 (6%)	2	2
3	D	95/176 (54%)	83 (87%)	6 (6%)	6 (6%)	2	1
All	All	683/804 (85%)	607 (89%)	50 (7%)	26 (4%)	4	5

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	GLY
3	C	137	GLU
3	C	167	VAL
3	C	168	SER
3	D	100	MET
1	A	138	ASP
1	A	177	ASN
3	C	92	LYS
3	C	165	ILE
3	C	166	THR
3	D	48	SER
3	D	65	LEU
3	D	67	ARG
1	A	231	GLU
1	A	245	SER
1	A	248	ILE
2	B	237	SER
3	C	55	GLN
3	C	94	LYS
3	C	159	GLU
3	C	160	HIS
3	D	49	MET

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Mol	Chain	Res	Type
3	D	95	SER
2	B	195	SER
2	B	278	GLY
1	A	157	GLY

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	184/195 (94%)	162 (88%)	22 (12%)	6	11
2	B	178/191 (93%)	160 (90%)	18 (10%)	9	16
3	C	159/161 (99%)	128 (80%)	31 (20%)	2	2
3	D	93/161 (58%)	78 (84%)	15 (16%)	3	4
All	All	614/708 (87%)	528 (86%)	86 (14%)	4	7

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	THR
1	A	61	ASN
1	A	112	LYS
1	A	118	LEU
1	A	138	ASP
1	A	146	ARG
1	A	150	LYS
1	A	172	VAL
1	A	175	GLU
1	A	176	LEU
1	A	177	ASN
1	A	178	MET
1	A	186	LEU
1	A	193	VAL
1	A	194	THR
1	A	202	LEU
1	A	206	THR

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Mol	Chain	Res	Type
1	A	213	LEU
1	A	245	SER
1	A	246	GLN
1	A	249	THR
1	A	251	GLU
2	B	174	ILE
2	B	184	LEU
2	B	212	ASP
2	B	228	ASN
2	B	260	ARG
2	B	267	THR
2	B	302	ARG
2	B	312	VAL
2	B	313	ILE
2	B	314	GLU
2	B	316	GLN
2	B	317	SER
2	B	319	LYS
2	B	344	LYS
2	B	354	LYS
2	B	356	ARG
2	B	375	LEU
2	B	384	THR
3	C	5	PHE
3	C	10	GLN
3	C	12	ARG
3	C	28	LYS
3	C	35	SER
3	C	42	ARG
3	C	49	MET
3	C	50	THR
3	C	57	SER
3	C	59	LEU
3	C	64	LYS
3	C	69	LEU
3	C	74	ILE
3	C	77	VAL
3	C	78	LEU
3	C	84	LYS
3	C	90	LEU
3	C	91	ASP
3	C	96	SER

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Mol	Chain	Res	Type
3	C	99	ILE
3	C	105	GLU
3	C	113	GLN
3	C	127	LEU
3	C	139	GLU
3	C	144	LEU
3	C	153	LEU
3	C	156	LEU
3	C	160	HIS
3	C	164	ILE
3	C	166	THR
3	C	169	ASP
3	D	5	PHE
3	D	10	GLN
3	D	26	ARG
3	D	28	LYS
3	D	38	LEU
3	D	52	MET
3	D	59	LEU
3	D	61	ASN
3	D	66	GLN
3	D	67	ARG
3	D	69	LEU
3	D	78	LEU
3	D	87	LEU
3	D	90	LEU
3	D	99	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	46	ASN
1	A	61	ASN
1	A	67	GLN
1	A	126	GLN
1	A	127	GLN
1	A	177	ASN
2	B	173	ASN
2	B	228	ASN
2	B	242	HIS
2	B	274	ASN

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Mol	Chain	Res	Type
2	B	289	ASN
2	B	316	GLN
3	C	10	GLN
3	C	22	ASN
3	C	44	HIS
3	C	113	GLN
3	D	10	GLN
3	D	55	GLN
3	D	61	ASN

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	219/234 (93%)	1.00	27 (12%) 5 3	39, 56, 120, 128	0
2	B	204/218 (93%)	1.17	31 (15%) 3 1	32, 66, 121, 130	0
3	C	173/176 (98%)	0.62	10 (5%) 26 20	40, 73, 99, 107	0
3	D	97/176 (55%)	1.60	33 (34%) 0 0	56, 93, 117, 119	0
All	All	693/804 (86%)	1.04	101 (14%) 3 2	32, 70, 119, 130	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	220	TYR	8.9
1	A	36	MET	7.4
2	B	174	ILE	7.3
3	D	89	TRP	7.0
1	A	34	ALA	6.6
1	A	57	GLU	6.3
3	D	48	SER	6.2
3	D	99	ILE	5.8
2	B	224	MET	5.8
3	D	44	HIS	5.4
3	D	97	PHE	5.0
3	D	50	THR	5.0
2	B	215	ILE	4.9
1	A	244	TYR	4.7
3	D	87	LEU	4.7
3	D	98	LEU	4.6
3	D	5	PHE	4.6
3	D	40	PHE	4.6
3	D	90	LEU	4.5
1	A	246	GLN	4.5
2	B	216	ARG	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	137	GLN	4.4
3	D	43	LEU	4.3
3	D	94	LYS	4.3
3	D	33	TRP	4.1
2	B	183	LYS	4.0
3	D	49	MET	3.9
1	A	47	LEU	3.9
3	D	93	SER	3.9
3	D	78	LEU	3.9
3	D	82	ARG	3.8
1	A	54	HIS	3.7
3	D	84	LYS	3.7
1	A	55	LYS	3.7
3	D	41	CYS	3.6
2	B	230	VAL	3.6
1	A	43	PHE	3.6
3	D	51	VAL	3.4
2	B	278	GLY	3.3
1	A	75	ILE	3.3
1	A	59	ARG	3.2
3	C	52	MET	3.2
2	B	185	MET	3.2
1	A	82	SER	3.1
2	B	181	LEU	3.0
3	D	96	SER	3.0
2	B	235	TYR	3.0
3	D	6	GLU	3.0
2	B	233	GLU	3.0
2	B	175	SER	2.9
2	B	199	ALA	2.9
3	D	46	GLN	2.8
1	A	56	GLN	2.8
1	A	68	PHE	2.8
3	D	100	MET	2.8
1	A	94	GLY	2.8
1	A	71	MET	2.8
1	A	93	VAL	2.7
3	D	86	ASN	2.7
3	D	60	PHE	2.7
2	B	285	GLU	2.7
1	A	46	ASN	2.7
3	D	37	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	64	PHE	2.6
1	A	81	ALA	2.5
2	B	225	GLY	2.5
3	C	120	GLN	2.5
1	A	133	GLY	2.5
2	B	276	ALA	2.4
3	C	165	ILE	2.4
2	B	338	THR	2.4
2	B	240	GLN	2.4
3	C	163	GLU	2.4
2	B	356	ARG	2.3
1	A	221	LEU	2.3
3	C	161	LYS	2.3
3	C	169	ASP	2.3
2	B	263	ILE	2.3
1	A	60	LYS	2.3
3	C	141	PHE	2.3
2	B	367	CYS	2.3
2	B	353	ALA	2.2
3	C	160	HIS	2.2
3	D	91	ASP	2.2
3	C	6	GLU	2.2
2	B	382	PHE	2.2
3	D	95	SER	2.1
1	A	247	GLU	2.1
2	B	323	MET	2.1
3	D	68	LYS	2.1
1	A	165	GLY	2.1
2	B	229	PRO	2.1
3	D	75	GLN	2.1
2	B	303	LEU	2.1
2	B	218	LYS	2.1
3	C	112	TYR	2.1
3	D	38	LEU	2.1
2	B	304	ARG	2.1
2	B	350	VAL	2.0
1	A	119	ILE	2.0
2	B	342	PHE	2.0

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

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.