



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:22 AM UTC

PDB ID : 1Q1J / pdb_00001q1j
Title : Crystal Structure Analysis of anti-HIV-1 Fab 447-52D in complex with V3 peptide
Authors : Stanfield, R.L.; Gorny, M.K.; Williams, C.; Zolla-Pazner, S.; Wilson, I.A.
Deposited on : 2003-07-21
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

<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

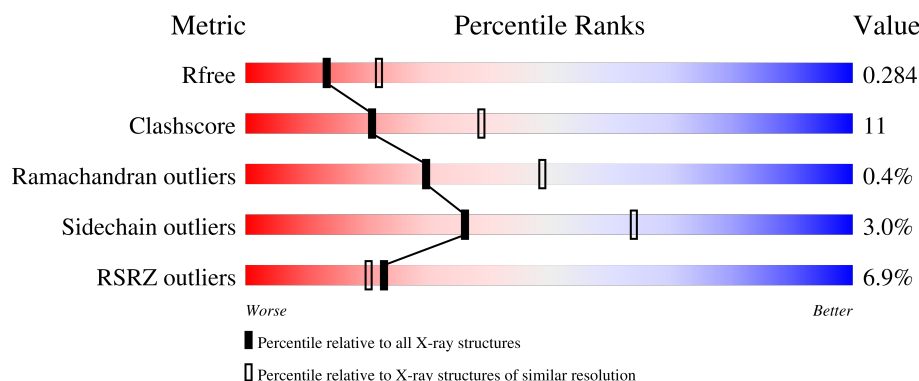
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 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	L	215	2% 72% 27% .
1	M	215	7% 71% 27% .
2	H	231	6% 74% 25% .
2	I	231	13% 73% 26% .
3	P	16	6% 62% 38% .

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Mol	Chain	Length	Quality of chain
3	Q	16	12%25%25%12%38%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6875 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 Fab 447-52D, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	215	Total	C	N	O	S	0	0	0
			1598	1004	265	325	4			
1	M	215	Total	C	N	O	S	0	0	0
			1598	1004	265	325	4			

- Molecule 2 is a protein called Fab 447-52D, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	231	Total	C	N	O	S	0	0	0
			1740	1095	291	346	8			
2	I	231	Total	C	N	O	S	0	0	0
			1740	1095	291	346	8			

- Molecule 3 is a protein called gp120 V3 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	10	Total	C	N	O	0	0	0
			73	45	18	10			
3	Q	10	Total	C	N	O	0	0	0
			73	45	18	10			

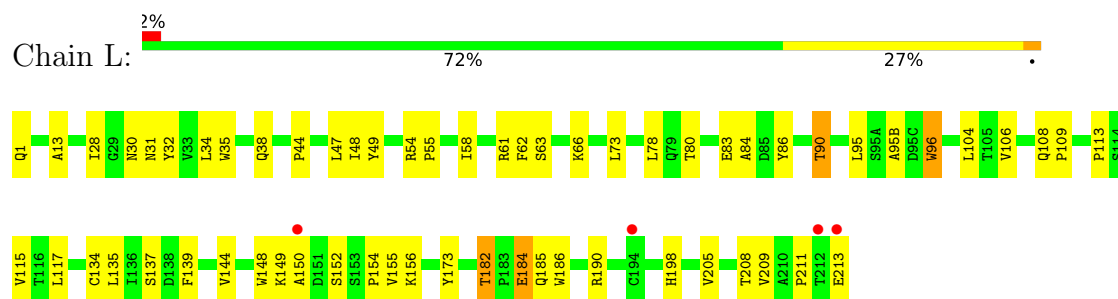
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	13	Total	O	0	0
			13	13		
4	H	16	Total	O	0	0
			16	16		
4	M	13	Total	O	0	0
			13	13		
4	I	11	Total	O	0	0
			11	11		

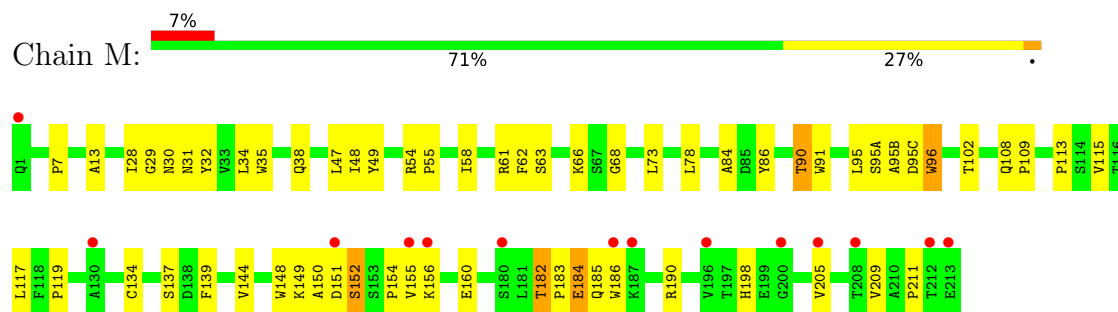
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

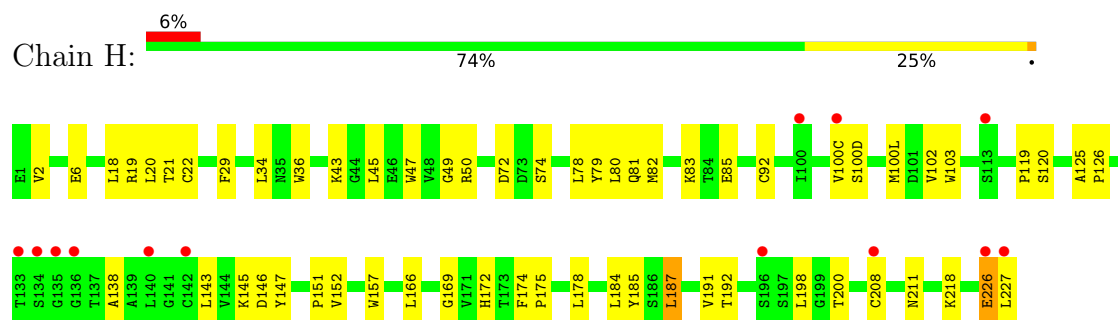
- Molecule 1: Fab 447-52D, light chain



- Molecule 1: Fab 447-52D, light chain

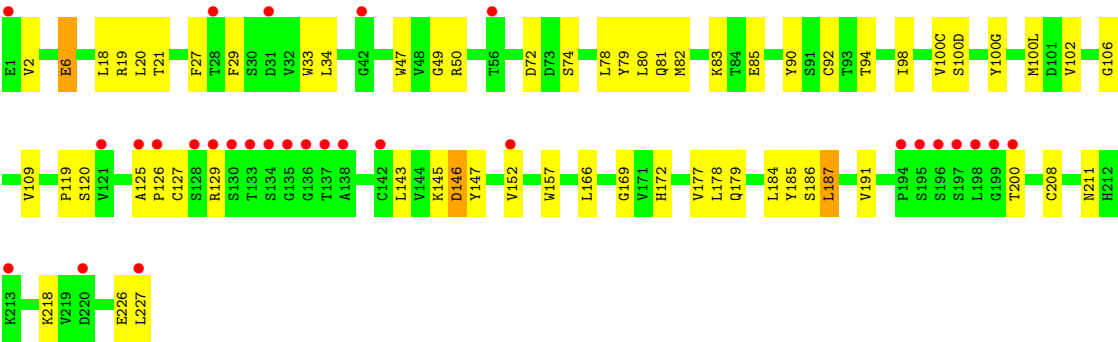


- Molecule 2: Fab 447-52D, heavy chain

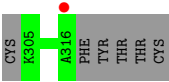


- Molecule 2: Fab 447-52D, heavy chain

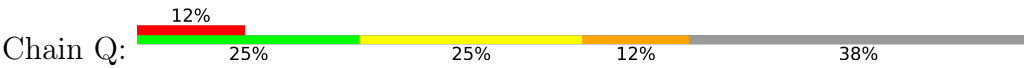




● Molecule 3: gp120 V3 peptide



● Molecule 3: gp120 V3 peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.67Å 74.90Å 100.05Å 90.00° 100.57° 90.00°	Depositor
Resolution (Å)	29.80 – 2.50 29.80 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.8 (29.80-2.50) 90.7 (29.80-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.250 , 0.285 0.249 , 0.284	Depositor DCC
R_{free} test set	2000 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	40.9	Xtriage
Anisotropy	0.396	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6875	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% 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	L	0.45	0/1639	0.94	7/2239 (0.3%)
1	M	0.44	0/1639	0.93	6/2239 (0.3%)
2	H	0.48	0/1779	0.91	2/2421 (0.1%)
2	I	0.45	0/1779	0.90	2/2421 (0.1%)
3	P	0.81	0/74	1.31	0/98
3	Q	1.82	3/74 (4.1%)	1.37	1/98 (1.0%)
All	All	0.50	3/6984 (0.0%)	0.93	18/9516 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	Q	307	ILE	N-CA	7.52	1.55	1.46
3	Q	313	PRO	CA-C	6.38	1.58	1.52
3	Q	306	ARG	CA-C	6.11	1.59	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Q	313	PRO	CB-CA-C	-8.21	102.61	111.56
1	L	96	TRP	N-CA-C	-7.46	98.41	110.20
1	M	96	TRP	N-CA-C	-6.77	98.99	109.76
1	L	31	ASN	N-CA-C	6.57	119.90	109.65
1	M	31	ASN	N-CA-C	6.54	119.86	109.65
1	M	182	THR	N-CA-C	-6.41	101.23	110.40
1	L	61	ARG	N-CA-C	-6.33	104.82	112.54
1	M	61	ARG	N-CA-C	-6.27	104.89	112.54
1	L	182	THR	N-CA-C	-6.26	101.44	110.40
2	I	120	SER	N-CA-C	-6.17	99.18	109.24
2	H	120	SER	N-CA-C	-5.99	99.48	109.24
1	L	95	LEU	N-CA-C	5.92	117.74	111.28
1	M	95	LEU	N-CA-C	5.74	117.53	111.28
1	L	137	SER	N-CA-C	5.34	117.10	109.14
1	M	137	SER	N-CA-C	5.23	116.93	109.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	173	TYR	N-CA-C	5.15	117.99	109.85
2	H	226	GLU	N-CA-C	5.06	116.89	109.24
2	I	98	ILE	N-CA-C	-5.03	101.13	108.17

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1598	0	1552	36	0
1	M	1598	0	1552	44	0
2	H	1740	0	1706	40	0
2	I	1740	0	1706	43	0
3	P	73	0	74	0	0
3	Q	73	0	74	11	0
4	H	16	0	0	1	0
4	I	11	0	0	0	0
4	L	13	0	0	1	0
4	M	13	0	0	0	0
All	All	6875	0	6664	149	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 (149) 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:I:100(G):TYR:CE2	3:Q:307:ILE:HD12	2.00	0.96
2:I:211:ASN:ND2	2:I:218:LYS:HE2	1.85	0.92
2:H:211:ASN:ND2	2:H:218:LYS:HE2	1.90	0.86
2:I:100(G):TYR:CZ	3:Q:307:ILE:HD12	2.14	0.82
1:M:95(B):ALA:HB3	2:I:50:ARG:NH2	1.96	0.81
2:H:20:LEU:HG	2:H:82:MET:HE2	1.63	0.80
2:I:211:ASN:HD21	2:I:218:LYS:HE2	1.47	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:95(B):ALA:HB3	2:H:50:ARG:NH2	1.99	0.77
2:I:20:LEU:HG	2:I:82:MET:HE2	1.65	0.77
1:M:32:TYR:O	1:M:90:THR:HG23	1.87	0.74
1:L:32:TYR:O	1:L:90:THR:HG23	1.88	0.74
2:H:2:VAL:HG11	2:H:102:VAL:HG21	1.70	0.73
2:I:100(G):TYR:CE2	3:Q:307:ILE:CD1	2.72	0.71
2:H:211:ASN:HD21	2:H:218:LYS:HE2	1.53	0.71
1:M:91:TRP:HB2	3:Q:313:PRO:HB3	1.73	0.69
2:I:72:ASP:OD1	2:I:74:SER:HB3	1.92	0.69
1:L:55:PRO:HD2	1:L:58:ILE:HG13	1.75	0.69
2:H:72:ASP:OD1	2:H:74:SER:HB3	1.94	0.67
2:H:143:LEU:HG	2:H:145:LYS:HG3	1.77	0.67
2:H:226:GLU:O	2:H:227:LEU:HB2	1.94	0.67
1:M:55:PRO:HD2	1:M:58:ILE:HG13	1.77	0.66
2:I:143:LEU:HG	2:I:145:LYS:HG3	1.78	0.66
2:I:226:GLU:O	2:I:227:LEU:HB2	1.96	0.65
2:I:125:ALA:HB1	2:I:227:LEU:HD12	1.80	0.64
2:I:82:MET:HE1	2:I:109:VAL:HG21	1.79	0.64
1:L:113:PRO:HB3	1:L:139:PHE:HB3	1.78	0.64
2:I:2:VAL:HG11	2:I:102:VAL:HG21	1.80	0.63
1:M:91:TRP:CB	3:Q:313:PRO:HB3	2.29	0.62
1:M:113:PRO:HB3	1:M:139:PHE:HB3	1.81	0.62
2:H:125:ALA:HB1	2:H:227:LEU:HD12	1.82	0.62
1:L:182:THR:OG1	1:L:184:GLU:HG3	1.99	0.62
2:H:198:LEU:HA	4:H:230:HOH:O	2.00	0.61
2:I:126:PRO:HG2	2:I:227:LEU:HD13	1.81	0.61
1:M:182:THR:OG1	1:M:184:GLU:HG3	2.01	0.60
2:I:83:LYS:HE2	2:I:85:GLU:CD	2.26	0.60
1:M:182:THR:OG1	1:M:185:GLN:HG3	2.02	0.59
2:H:126:PRO:HG2	2:H:227:LEU:HD13	1.85	0.59
1:L:182:THR:HG1	1:L:185:GLN:HG3	1.67	0.59
2:H:83:LYS:HE2	2:H:85:GLU:CD	2.27	0.58
1:L:182:THR:OG1	1:L:185:GLN:HG3	2.03	0.58
2:H:100(C):VAL:HG13	2:H:100(D):SER:N	2.19	0.58
1:M:34:LEU:HD23	1:M:49:TYR:HA	1.86	0.58
1:L:54:ARG:HD3	1:L:62:PHE:O	2.02	0.58
2:H:187:LEU:C	2:H:187:LEU:HD12	2.29	0.57
1:M:54:ARG:HD3	1:M:62:PHE:O	2.04	0.57
1:L:34:LEU:HD23	1:L:49:TYR:HA	1.87	0.57
1:M:38:GLN:O	1:M:84:ALA:HB1	2.04	0.57
1:L:95(B):ALA:HB3	2:H:50:ARG:HH22	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:38:GLN:O	1:L:84:ALA:HB1	2.06	0.56
1:M:96:TRP:CD1	2:I:100(L):MET:HE3	2.40	0.56
2:I:100(C):VAL:HG13	2:I:100(D):SER:N	2.22	0.55
2:I:187:LEU:C	2:I:187:LEU:HD12	2.31	0.55
1:M:160:GLU:OE1	2:I:179:GLN:HG2	2.06	0.55
1:M:95(B):ALA:HB3	2:I:50:ARG:HH22	1.68	0.55
2:I:119:PRO:HB3	2:I:147:TYR:HB3	1.89	0.54
2:H:119:PRO:HB3	2:H:147:TYR:HB3	1.90	0.54
1:L:213:GLU:OXT	1:L:213:GLU:HG2	2.07	0.54
2:I:6:GLU:CD	2:I:106:GLY:H	2.16	0.53
2:H:21:THR:HG22	2:H:79:TYR:CD1	2.44	0.53
1:L:96:TRP:CD1	2:H:100(L):MET:HE3	2.44	0.53
2:I:21:THR:HG22	2:I:79:TYR:CD1	2.45	0.52
2:I:169:GLY:O	2:I:191:VAL:HA	2.10	0.52
1:L:186:TRP:CE2	1:L:211:PRO:HG3	2.44	0.51
2:H:169:GLY:O	2:H:191:VAL:HA	2.09	0.51
2:H:45:LEU:HD11	2:H:103:TRP:HZ3	1.76	0.51
1:M:186:TRP:CE2	1:M:211:PRO:HG3	2.47	0.50
1:M:91:TRP:HB3	3:Q:313:PRO:HG3	1.92	0.50
1:M:182:THR:HG1	1:M:185:GLN:HG3	1.76	0.50
2:I:178:LEU:HD13	2:I:185:TYR:CE2	2.47	0.50
2:I:33:TRP:HH2	3:Q:314:GLY:O	1.93	0.50
1:M:7:PRO:O	1:M:102:THR:OG1	2.22	0.49
1:M:149:LYS:HD3	1:M:154:PRO:HA	1.94	0.49
2:I:34:LEU:HD13	2:I:78:LEU:HD13	1.94	0.49
2:I:157:TRP:CH2	2:I:208:CYS:HB3	2.48	0.49
2:I:18:LEU:HB3	2:I:82:MET:HE3	1.95	0.49
2:H:19:ARG:HD2	2:H:81:GLN:OE1	2.12	0.49
1:L:1:GLN:OE1	2:H:43:LYS:HA	2.12	0.48
1:L:149:LYS:HD3	1:L:154:PRO:HA	1.95	0.48
1:L:83:GLU:HG3	1:L:104:LEU:O	2.13	0.48
4:L:224:HOH:O	2:H:175:PRO:HD2	2.12	0.48
2:H:157:TRP:CH2	2:H:208:CYS:HB3	2.48	0.48
2:H:34:LEU:HD13	2:H:78:LEU:HD13	1.95	0.48
2:I:178:LEU:HD13	2:I:185:TYR:CZ	2.48	0.48
1:M:117:LEU:HD21	1:M:209:VAL:HG22	1.95	0.48
2:I:19:ARG:HD2	2:I:81:GLN:OE1	2.12	0.48
1:M:13:ALA:HB3	1:M:78:LEU:CD1	2.44	0.48
2:I:146:ASP:HB3	2:I:184:LEU:HD13	1.96	0.48
1:L:13:ALA:HB3	1:L:78:LEU:CD1	2.43	0.48
2:H:47:TRP:CZ2	2:H:49:GLY:HA2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:117:LEU:HD21	1:L:209:VAL:HG22	1.95	0.47
1:L:186:TRP:O	1:L:211:PRO:HG2	2.14	0.47
2:H:18:LEU:HB3	2:H:82:MET:HE3	1.96	0.47
1:M:186:TRP:O	1:M:211:PRO:HG2	2.15	0.47
2:I:47:TRP:CZ2	2:I:49:GLY:HA2	2.49	0.47
2:I:177:VAL:HG12	2:I:186:SER:O	2.15	0.47
1:M:91:TRP:CE3	3:Q:313:PRO:HD3	2.51	0.46
1:M:108:GLN:HB2	1:M:109:PRO:HD2	1.98	0.46
1:L:47:LEU:HD11	1:L:86:TYR:HE1	1.81	0.46
2:I:29:PHE:CZ	2:I:34:LEU:HD11	2.51	0.46
2:H:178:LEU:HD13	2:H:185:TYR:CZ	2.50	0.46
1:M:29:GLY:HA2	1:M:68:GLY:O	2.16	0.45
1:L:134:CYS:HB2	1:L:148:TRP:CH2	2.51	0.45
1:L:80:THR:HA	1:L:106:VAL:HG11	1.99	0.45
2:I:90:TYR:O	2:I:106:GLY:HA2	2.16	0.45
2:H:2:VAL:CG1	2:H:102:VAL:HG21	2.45	0.45
1:L:108:GLN:HB2	1:L:109:PRO:HD2	1.99	0.45
1:L:115:VAL:HG21	1:L:205:VAL:HG21	1.98	0.45
1:M:150:ALA:O	1:M:152:SER:N	2.48	0.45
2:H:146:ASP:HB3	2:H:184:LEU:HD13	1.99	0.44
2:H:29:PHE:CZ	2:H:34:LEU:HD11	2.53	0.44
1:M:150:ALA:O	1:M:151:ASP:HB2	2.18	0.44
1:M:134:CYS:HB2	1:M:148:TRP:CH2	2.54	0.43
1:M:32:TYR:CD1	3:Q:309:ILE:CG2	3.02	0.43
2:H:100(C):VAL:CG1	2:H:100(D):SER:N	2.81	0.43
1:M:91:TRP:HA	1:M:95(C):ASP:O	2.19	0.43
2:I:33:TRP:CE2	3:Q:315:ARG:NH1	2.87	0.43
1:M:47:LEU:HD11	1:M:86:TYR:HE1	1.84	0.43
1:L:144:VAL:HG12	1:L:198:HIS:HB2	2.01	0.43
1:M:182:THR:O	1:M:183:PRO:C	2.62	0.42
2:I:34:LEU:HD23	2:I:94:THR:HG22	2.01	0.42
1:L:63:SER:O	1:L:73:LEU:HD12	2.19	0.42
1:L:208:THR:HG22	1:L:209:VAL:N	2.34	0.42
1:M:28:ILE:O	1:M:66:LYS:HE3	2.20	0.42
1:M:95(A):SER:O	2:I:50:ARG:NH2	2.53	0.42
1:M:144:VAL:HG12	1:M:198:HIS:HB2	2.00	0.42
2:H:22:CYS:HB3	2:H:78:LEU:HB3	2.02	0.42
1:L:95(B):ALA:CB	2:H:50:ARG:HH22	2.33	0.41
1:M:95(B):ALA:CB	2:I:50:ARG:HH22	2.32	0.41
2:I:20:LEU:HD12	2:I:80:LEU:HD23	2.02	0.41
1:L:28:ILE:O	1:L:66:LYS:HE3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:44:PRO:HD2	2:H:103:TRP:CG	2.56	0.41
2:H:227:LEU:C	2:H:227:LEU:HD23	2.45	0.41
1:M:32:TYR:CD1	3:Q:309:ILE:HG23	2.55	0.41
1:M:182:THR:CB	1:M:184:GLU:HG3	2.51	0.41
2:I:2:VAL:HG13	2:I:27:PHE:CD1	2.55	0.41
1:M:63:SER:O	1:M:73:LEU:HD12	2.20	0.41
2:H:178:LEU:HD13	2:H:185:TYR:CE2	2.56	0.41
1:L:35:TRP:HB2	1:L:48:ILE:HB	2.03	0.41
1:L:190:ARG:HG3	1:L:190:ARG:NH1	2.36	0.41
1:L:135:LEU:HB3	2:H:174:PHE:CZ	2.56	0.41
2:H:36:TRP:CE2	2:H:80:LEU:HB2	2.56	0.41
1:M:119:PRO:HD3	2:I:127:CYS:HB2	2.03	0.41
2:H:138:ALA:O	2:H:192:THR:HA	2.22	0.40
1:M:35:TRP:HB2	1:M:48:ILE:HB	2.02	0.40
1:M:150:ALA:HB2	1:M:155:VAL:HG21	2.04	0.40
1:L:150:ALA:HB2	1:L:155:VAL:HG21	2.04	0.40
1:M:190:ARG:HH11	1:M:190:ARG:HG3	1.87	0.40
1:L:182:THR:CB	1:L:184:GLU:HG3	2.50	0.40
1:M:115:VAL:HG21	1:M:205:VAL:HG21	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	L	213/215 (99%)	202 (95%)	10 (5%)	1 (0%)	24	43
1	M	213/215 (99%)	201 (94%)	11 (5%)	1 (0%)	24	43
2	H	229/231 (99%)	216 (94%)	13 (6%)	0	100	100
2	I	229/231 (99%)	214 (93%)	13 (6%)	2 (1%)	14	27
3	P	8/16 (50%)	7 (88%)	1 (12%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	Q	8/16 (50%)	8 (100%)	0	0	100	100
All	All	900/924 (97%)	848 (94%)	48 (5%)	4 (0%)	30	49

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	M	152	SER
2	I	129	ARG
2	I	146	ASP
1	L	152	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	L	179/179 (100%)	175 (98%)	4 (2%)	45	73
1	M	179/179 (100%)	175 (98%)	4 (2%)	45	73
2	H	196/196 (100%)	188 (96%)	8 (4%)	27	53
2	I	196/196 (100%)	189 (96%)	7 (4%)	31	58
3	P	6/13 (46%)	6 (100%)	0	100	100
3	Q	6/13 (46%)	6 (100%)	0	100	100
All	All	762/776 (98%)	739 (97%)	23 (3%)	36	64

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	30	ASN
1	L	90	THR
1	L	156	LYS
1	L	184	GLU
2	H	6	GLU
2	H	92	CYS
2	H	151	PRO

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Mol	Chain	Res	Type
2	H	152	VAL
2	H	166	LEU
2	H	172	HIS
2	H	187	LEU
2	H	200	THR
1	M	30	ASN
1	M	90	THR
1	M	156	LYS
1	M	184	GLU
2	I	6	GLU
2	I	92	CYS
2	I	152	VAL
2	I	166	LEU
2	I	172	HIS
2	I	187	LEU
2	I	200	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	198	HIS
2	H	179	GLN
2	H	211	ASN
2	I	179	GLN
2	I	211	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	L	215/215 (100%)	0.40	4 (1%) 66 62	27, 43, 60, 77	0
1	M	215/215 (100%)	0.73	14 (6%) 25 22	29, 48, 78, 88	0
2	H	231/231 (100%)	0.56	13 (5%) 30 26	20, 40, 65, 92	0
2	I	231/231 (100%)	0.81	29 (12%) 8 6	29, 44, 74, 103	0
3	P	10/16 (62%)	1.29	1 (10%) 12 10	45, 51, 67, 68	0
3	Q	10/16 (62%)	1.72	2 (20%) 3 2	49, 54, 66, 67	0
All	All	912/924 (98%)	0.65	63 (6%) 23 20	20, 44, 72, 103	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	198	LEU	6.4
2	H	227	LEU	5.8
2	I	135	GLY	5.0
2	H	135	GLY	4.8
2	H	142	CYS	4.6
2	I	130	SER	4.6
2	I	134	SER	4.3
3	Q	316	ALA	4.2
3	P	316	ALA	4.2
2	I	137	THR	4.1
2	I	227	LEU	3.9
2	I	133	THR	3.9
2	I	136	GLY	3.8
2	H	133	THR	3.8
2	I	197	SER	3.7
2	I	200	THR	3.7
2	H	134	SER	3.5
2	I	128	SER	3.5
1	M	1	GLN	3.2

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Mol	Chain	Res	Type	RSRZ
2	I	199	GLY	3.2
2	I	1	GLU	3.1
2	H	100(C)	VAL	3.0
1	L	212	THR	3.0
2	I	129	ARG	2.7
3	Q	308	HIS	2.7
2	I	142	CYS	2.6
1	M	205	VAL	2.6
2	I	126	PRO	2.6
2	H	196	SER	2.6
2	I	195	SER	2.6
1	L	194	CYS	2.6
2	I	220	ASP	2.5
2	H	140	LEU	2.5
1	M	130	ALA	2.4
1	M	212	THR	2.3
1	M	186	TRP	2.3
2	H	113	SER	2.3
2	I	194	PRO	2.3
2	I	196	SER	2.3
1	L	213	GLU	2.3
1	M	155	VAL	2.3
1	M	156	LYS	2.3
2	H	208	CYS	2.3
1	M	200	GLY	2.2
1	M	187	LYS	2.2
1	M	180	SER	2.2
2	H	226	GLU	2.2
2	H	100	ILE	2.2
1	M	196	VAL	2.1
2	I	152	VAL	2.1
1	L	150	ALA	2.1
1	M	151	ASP	2.1
2	I	121	VAL	2.1
2	I	138	ALA	2.1
2	H	136	GLY	2.1
2	I	42	GLY	2.1
2	I	28	THR	2.1
2	I	125	ALA	2.0
1	M	208	THR	2.0
2	I	213	LYS	2.0
2	I	31	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	M	213	GLU	2.0
2	I	56	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.