



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:36 AM UTC

PDB ID : 1SVZ / pdb_00001svz
Title : Crystal structure of the single-chain Fv fragment 1696 in complex with the epitope peptide corresponding to N-terminus of HIV-2 protease
Authors : Rezacova, P.; Brynda, J.; Lescar, J.; Bentley, G.A.; Fabry, M.; Horejsi, M.; Sedlacek, J.
Deposited on : 2004-03-30
Resolution : 1.89 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

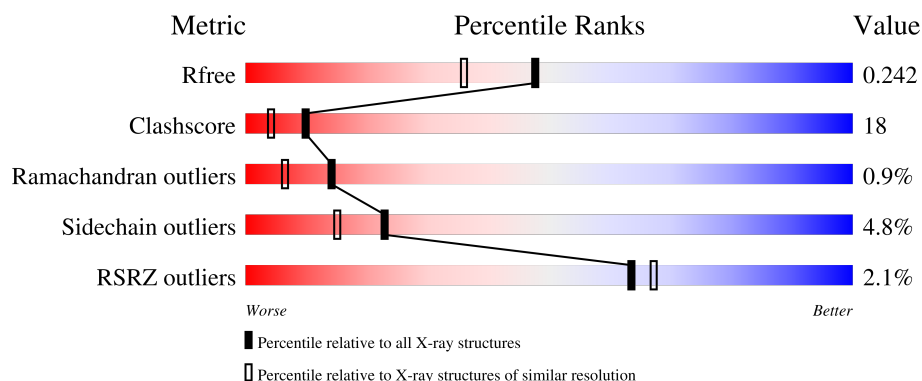
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	247	
1	B	247	
2	C	8	
2	D	8	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4146 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 single-chain Fv fragment 1696.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	2	4	0
			1847	1176	301	364	6			
1	B	232	Total	C	N	O	S	0	3	0
			1834	1164	300	364	6			

- Molecule 2 is a protein called epitope peptide corresponding to N-terminus of HIV-2 protease.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	0	0	0
			55	39	8	8			
2	D	6	Total	C	N	O	0	0	0
			55	39	8	8			

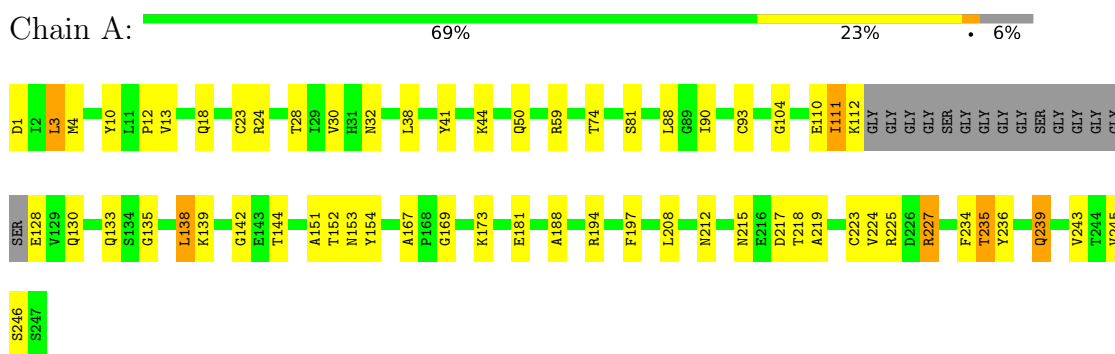
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	202	Total	O	0	0
			202	202		
3	B	137	Total	O	0	0
			137	137		
3	C	13	Total	O	0	0
			13	13		
3	D	3	Total	O	0	0
			3	3		

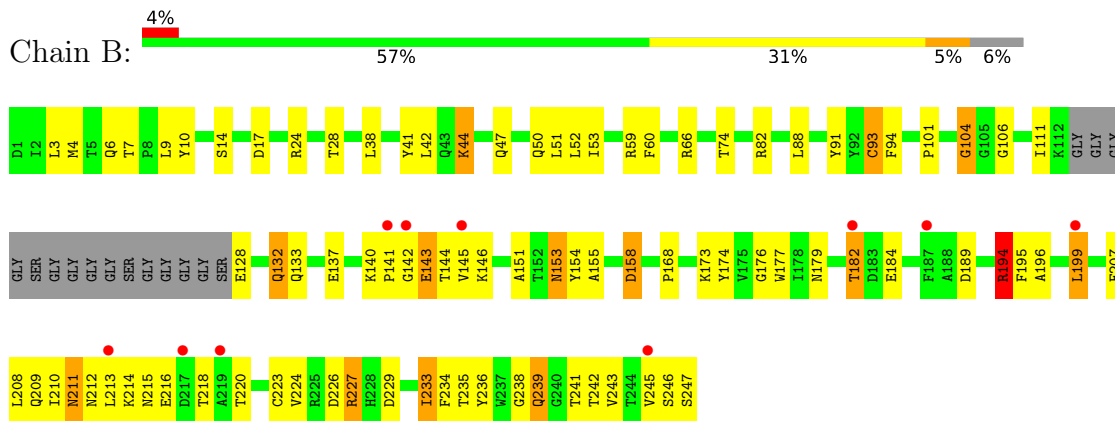
3 Residue-property plots

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.

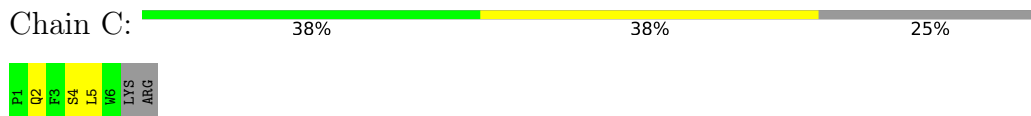
- Molecule 1: single-chain Fv fragment 1696



- Molecule 1: single-chain Fv fragment 1696



- Molecule 2: epitope peptide corresponding to N-terminus of HIV-2 protease



- Molecule 2: epitope peptide corresponding to N-terminus of HIV-2 protease



P1	Q2	F3	S4	L5	W6	LYS	ARG
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	44.18Å 56.41Å 88.30Å 90.00° 102.13° 90.00°	Depositor
Resolution (Å)	19.79 – 1.89 19.79 – 1.89	Depositor EDS
% Data completeness (in resolution range)	95.8 (19.79-1.89) 97.2 (19.79-1.89)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 1.89Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.196 , 0.239 0.199 , 0.242	Depositor DCC
R_{free} test set	1621 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 67.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4146	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.86% 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	0.67	0/1913	1.06	12/2598 (0.5%)
1	B	0.57	0/1896	1.01	7/2573 (0.3%)
2	C	0.54	0/58	1.29	2/78 (2.6%)
2	D	0.46	0/58	1.05	0/78
All	All	0.62	0/3925	1.04	21/5327 (0.4%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	135	GLY	N-CA-C	8.88	123.02	112.10
1	B	227	ARG	N-CA-C	-7.46	101.28	110.61
1	A	227	ARG	N-CA-C	-7.30	99.62	110.06
1	A	223	CYS	N-CA-C	-7.20	98.58	110.17
1	A	104	GLY	N-CA-C	-6.96	103.22	112.82
1	B	104	GLY	N-CA-C	-6.73	102.22	112.51
1	B	153	ASN	N-CA-C	6.33	120.31	112.58
1	B	223	CYS	N-CA-C	-6.27	100.35	110.32
1	A	235[A]	THR	N-CA-C	6.21	119.33	111.69
1	A	235[B]	THR	N-CA-C	6.21	119.33	111.69
1	B	28	THR	N-CA-C	-5.98	102.06	110.50
1	A	153	ASN	N-CA-C	5.93	121.81	113.56
2	C	5	LEU	N-CA-C	5.76	120.15	113.18
1	B	238	GLY	N-CA-C	-5.67	103.72	112.85
1	A	93	CYS	N-CA-C	-5.64	101.68	110.42
1	A	188	ALA	N-CA-C	-5.46	101.57	110.20
1	B	93	CYS	N-CA-C	-5.28	102.65	110.46
1	A	30	VAL	N-CA-C	-5.26	102.06	109.21
1	A	167	ALA	CA-C-N	5.20	125.25	119.32
1	A	167	ALA	C-N-CA	5.20	125.25	119.32
2	C	4	SER	N-CA-C	5.11	120.14	112.94

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	A	1847	0	1757	52	0
1	B	1834	0	1749	78	0
2	C	55	0	52	2	0
2	D	55	0	52	3	0
3	A	202	0	0	12	0
3	B	137	0	0	12	0
3	C	13	0	0	0	0
3	D	3	0	0	0	0
All	All	4146	0	3610	130	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 18.

All (130) 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:133:GLN:H	1:B:239:GLN:HE22	1.02	0.99
1:A:130:GLN:HB3	1:A:152[B]:THR:CG2	1.94	0.96
1:A:142:GLY:HA2	1:A:212:ASN:HD22	1.31	0.96
1:B:214:LYS:HB2	3:B:327:HOH:O	1.67	0.94
1:A:227:ARG:HG3	1:A:235[A]:THR:HG21	1.47	0.93
1:A:133:GLN:H	1:A:239:GLN:HE22	1.17	0.92
1:B:133:GLN:H	1:B:239:GLN:NE2	1.72	0.88
1:A:239:GLN:HE21	1:A:239:GLN:H	1.22	0.87
1:A:142:GLY:HA2	1:A:212:ASN:ND2	1.91	0.85
1:A:111:ILE:H	1:A:111:ILE:HD13	1.44	0.82
1:B:179:ASN:HB3	1:B:182:THR:HG22	1.62	0.82
1:B:9:LEU:HD22	1:B:10:TYR:CE1	2.16	0.81
1:B:227:ARG:HB2	1:B:235[A]:THR:HG21	1.64	0.80
1:B:194:ARG:NH2	1:B:214:LYS:HG3	1.98	0.77
1:B:142:GLY:HA2	1:B:212:ASN:ND2	2.00	0.77
1:B:168:PRO:HA	3:B:249:HOH:O	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:GLN:N	1:B:239:GLN:HE22	1.83	0.76
1:A:90:ILE:HD12	3:A:362:HOH:O	1.84	0.75
1:A:215:ASN:HD22	1:B:215:ASN:HB2	1.53	0.73
1:B:42:LEU:O	1:B:50:GLN:HG2	1.89	0.73
1:B:142:GLY:HA2	1:B:212:ASN:HD22	1.52	0.72
1:B:140:LYS:HG2	1:B:246:SER:HB3	1.73	0.71
1:B:7:THR:HG23	3:B:296:HOH:O	1.90	0.70
1:A:130:GLN:HB3	1:A:152[B]:THR:HG23	1.73	0.68
3:B:306:HOH:O	2:D:1:PRO:HD3	1.94	0.68
1:B:218:THR:HA	1:B:243:VAL:O	1.92	0.68
1:A:1:ASP:HB2	3:A:384:HOH:O	1.95	0.66
1:B:220:THR:OG1	1:B:242:THR:HG22	1.97	0.65
1:B:128:GLU:HG3	1:B:153:ASN:ND2	2.12	0.64
1:A:181:GLU:HG3	2:C:2:GLN:NE2	2.13	0.63
1:B:145:VAL:HG12	1:B:213:LEU:HD21	1.80	0.62
1:B:4:MET:HB2	1:B:104:GLY:HA2	1.80	0.62
1:B:142:GLY:O	1:B:212:ASN:HA	2.01	0.61
1:B:179:ASN:CB	1:B:182:THR:HG22	2.31	0.61
1:B:140:LYS:O	1:B:143:GLU:HB2	2.00	0.60
1:A:197:PHE:CE1	1:A:208:LEU:HD13	2.37	0.60
1:B:235[A]:THR:HG23	1:B:236:TYR:CD2	2.37	0.59
1:B:14:SER:HB2	1:B:17:ASP:OD2	2.01	0.59
1:A:130:GLN:HB3	1:A:152[B]:THR:HG21	1.82	0.59
1:A:239:GLN:HE21	1:A:239:GLN:N	1.98	0.58
1:B:239:GLN:HE21	1:B:239:GLN:H	1.50	0.57
1:A:151:ALA:HB1	1:A:154:TYR:CE1	2.39	0.57
1:A:1:ASP:HB2	3:A:329:HOH:O	2.05	0.56
1:B:60:PHE:CE1	1:B:235[B]:THR:HG23	2.41	0.56
1:B:194:ARG:HG3	1:B:194:ARG:HH11	1.71	0.56
1:A:18:GLN:HG2	1:A:81:SER:O	2.05	0.56
1:B:199:LEU:HD23	1:B:199:LEU:N	2.21	0.55
1:A:181:GLU:HG3	2:C:2:GLN:HE21	1.70	0.55
1:A:130:GLN:O	1:A:152[B]:THR:HG22	2.06	0.55
1:B:9:LEU:HD22	1:B:10:TYR:CD1	2.42	0.55
1:B:226:ASP:HA	1:B:233:ILE:O	2.05	0.55
1:B:128:GLU:HA	3:B:308:HOH:O	2.08	0.54
1:A:139:LYS:HE2	1:A:144:THR:O	2.08	0.53
1:B:182:THR:HG21	3:B:348:HOH:O	2.08	0.53
1:A:13:VAL:O	1:A:111:ILE:HA	2.08	0.53
1:A:197:PHE:HE1	1:A:208:LEU:HD13	1.74	0.52
1:B:141:PRO:HB3	3:B:324:HOH:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:MET:HE3	1:A:23:CYS:SG	2.50	0.51
1:B:194:ARG:HB3	1:B:211:ASN:ND2	2.26	0.50
1:A:218:THR:O	1:A:219:ALA:HB2	2.11	0.50
1:A:235[A]:THR:HG23	1:A:236:TYR:CD2	2.46	0.50
1:A:169:GLY:N	1:B:242:THR:HG21	2.26	0.50
1:B:151:ALA:HB1	1:B:154:TYR:CE1	2.46	0.50
1:A:133:GLN:H	1:A:239:GLN:NE2	1.98	0.49
1:A:41:TYR:OH	1:A:234:PHE:HB2	2.12	0.49
1:A:173:LYS:HB3	1:A:173:LYS:NZ	2.28	0.49
1:B:141:PRO:HA	1:B:213:LEU:O	2.12	0.49
1:B:196:ALA:HB3	1:B:209:GLN:HB3	1.95	0.49
1:A:215:ASN:ND2	1:B:215:ASN:HB2	2.22	0.49
1:B:141:PRO:HB3	3:B:310:HOH:O	2.12	0.49
1:B:59:ARG:HH11	1:B:59:ARG:HG3	1.78	0.49
1:A:128:GLU:HG3	3:A:337:HOH:O	2.13	0.48
1:B:140:LYS:CG	1:B:246:SER:HB3	2.40	0.48
1:A:128:GLU:CG	3:A:337:HOH:O	2.61	0.48
1:B:41:TYR:OH	1:B:234:PHE:HB2	2.14	0.48
1:A:111:ILE:H	1:A:111:ILE:CD1	2.20	0.47
1:B:155:ALA:O	1:B:158:ASP:HB2	2.14	0.47
1:B:241:THR:HG21	3:B:311:HOH:O	2.14	0.47
1:B:141:PRO:HG2	1:B:247:SER:HA	1.95	0.47
1:A:24:ARG:HA	1:A:74:THR:O	2.15	0.47
1:A:32:ASN:ND2	3:A:313:HOH:O	2.47	0.47
1:B:59:ARG:HG3	1:B:59:ARG:NH1	2.30	0.47
1:B:174:TYR:CE2	1:B:176:GLY:HA2	2.50	0.47
1:B:239:GLN:NE2	1:B:239:GLN:H	2.14	0.46
1:A:111:ILE:HD13	1:A:111:ILE:N	2.23	0.46
1:A:194:ARG:HH22	1:A:217:ASP:CG	2.23	0.45
1:B:146:LYS:HE2	1:B:207:PHE:CD2	2.52	0.45
2:D:3:PHE:CE2	2:D:5:LEU:HB2	2.51	0.45
1:B:91:TYR:O	1:B:106:GLY:HA2	2.17	0.44
1:B:173:LYS:HE2	1:B:173:LYS:HB3	1.85	0.44
1:A:18:GLN:NE2	3:A:287:HOH:O	2.49	0.44
1:B:145:VAL:HG12	1:B:210:ILE:HB	2.00	0.44
1:B:132:GLN:HB2	3:B:268:HOH:O	2.15	0.44
1:B:137:GLU:HG2	3:B:311:HOH:O	2.16	0.44
1:A:112:LYS:HB2	3:A:416:HOH:O	2.18	0.44
1:B:44:LYS:HB2	1:B:47:GLN:NE2	2.33	0.44
1:A:128:GLU:N	3:A:371:HOH:O	2.51	0.43
1:B:194:ARG:HG3	1:B:194:ARG:NH1	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:3:PHE:HE2	2:D:6:TRP:CD2	2.36	0.43
1:A:88:LEU:HD11	1:A:111:ILE:HG23	2.00	0.43
1:B:53:ILE:HD13	1:B:59:ARG:HA	2.01	0.43
1:A:139:LYS:O	1:A:245:VAL:HA	2.18	0.43
1:A:218:THR:HA	1:A:243:VAL:O	2.19	0.43
1:B:151:ALA:HB1	1:B:154:TYR:HE1	1.84	0.43
1:B:182:THR:HG23	1:B:184:GLU:H	1.84	0.43
1:B:42:LEU:HB2	1:B:52:LEU:HD11	2.01	0.43
1:A:44:LYS:HE3	1:A:50:GLN:NE2	2.34	0.43
1:A:59:ARG:NH1	3:A:272:HOH:O	2.29	0.43
1:B:142:GLY:C	1:B:212:ASN:HA	2.42	0.42
1:A:128:GLU:CA	3:A:371:HOH:O	2.68	0.42
1:B:199:LEU:N	1:B:199:LEU:CD2	2.82	0.42
1:B:24:ARG:HA	1:B:74:THR:O	2.19	0.42
1:B:144:THR:OG1	1:B:211:ASN:HA	2.19	0.42
1:A:138:LEU:HD21	1:A:246:SER:HB2	2.02	0.42
1:B:94:PHE:CE2	1:B:101:PRO:HB2	2.55	0.41
1:B:88:LEU:HD11	1:B:111:ILE:HG23	2.02	0.41
1:B:224:VAL:HG11	1:B:234:PHE:HB3	2.02	0.41
1:B:141:PRO:CG	1:B:247:SER:HA	2.51	0.41
1:B:174:TYR:OH	1:B:177:TRP:HB3	2.21	0.41
1:B:229:ASP:OD1	1:B:229:ASP:N	2.54	0.41
1:A:128:GLU:OE1	1:A:128:GLU:HA	2.20	0.41
1:B:128:GLU:N	3:B:305:HOH:O	2.54	0.41
1:A:3:LEU:HD21	3:A:283:HOH:O	2.21	0.40
1:A:224:VAL:HG11	1:A:234:PHE:HB3	2.03	0.40
1:B:218:THR:OG1	1:B:245:VAL:HG23	2.22	0.40
1:B:6:GLN:HG2	1:B:93:CYS:SG	2.60	0.40
1:A:10:TYR:CD2	1:A:12:PRO:HG3	2.56	0.40
1:B:66:ARG:HD2	1:B:82:ARG:O	2.21	0.40
1:B:195:PHE:HD2	1:B:208:LEU:HD11	1.85	0.40
1:B:4:MET:CB	1:B:104:GLY:HA2	2.47	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	232/247 (94%)	223 (96%)	9 (4%)	0	100	100
1	B	231/247 (94%)	212 (92%)	15 (6%)	4 (2%)	7	2
2	C	4/8 (50%)	4 (100%)	0	0	100	100
2	D	4/8 (50%)	4 (100%)	0	0	100	100
All	All	471/510 (92%)	443 (94%)	24 (5%)	4 (1%)	14	8

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	194	ARG
1	B	189	ASP
1	B	143	GLU
1	B	233	ILE

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	206/205 (100%)	198 (96%)	8 (4%)	28	21
1	B	205/205 (100%)	193 (94%)	12 (6%)	18	10
2	C	6/8 (75%)	6 (100%)	0	100	100
2	D	6/8 (75%)	6 (100%)	0	100	100
All	All	423/426 (99%)	403 (95%)	20 (5%)	23	15

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	28	THR

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Mol	Chain	Res	Type
1	A	38	LEU
1	A	110	GLU
1	A	111	ILE
1	A	138	LEU
1	A	225	ARG
1	A	239	GLN
1	B	3	LEU
1	B	38	LEU
1	B	44	LYS
1	B	51	LEU
1	B	132	GLN
1	B	158	ASP
1	B	182	THR
1	B	194	ARG
1	B	199	LEU
1	B	211	ASN
1	B	216	GLU
1	B	239	GLN

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

Mol	Chain	Res	Type
1	A	18	GLN
1	A	35	ASN
1	A	50	GLN
1	A	98	HIS
1	A	132	GLN
1	A	212	ASN
1	A	215	ASN
1	A	239	GLN
1	B	18	GLN
1	B	27	GLN
1	B	32	ASN
1	B	35	ASN
1	B	47	GLN
1	B	58	ASN
1	B	130	GLN
1	B	153	ASN
1	B	209	GLN
1	B	211	ASN
1	B	212	ASN
1	B	239	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	232/247 (93%)	-0.21	0 100 100	17, 30, 54, 69	5 (2%)
1	B	232/247 (93%)	0.44	10 (4%) 40 42	25, 46, 69, 70	3 (1%)
2	C	6/8 (75%)	-0.14	0 100 100	25, 33, 38, 38	0
2	D	6/8 (75%)	0.76	0 100 100	54, 56, 62, 65	0
All	All	476/510 (93%)	0.12	10 (2%) 63 67	17, 38, 67, 70	8 (1%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	142	GLY	3.4
1	B	245	VAL	3.0
1	B	213	LEU	2.9
1	B	187	PHE	2.4
1	B	199	LEU	2.3
1	B	141	PRO	2.2
1	B	219	ALA	2.2
1	B	217	ASP	2.2
1	B	145	VAL	2.1
1	B	182	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.