



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:37 AM UTC

PDB ID : 3IDJ / pdb_00003idj
Title : Crystal structure of the HIV-1 Cross Neutralizing Monoclonal Antibody 2F5 Fab' fragment in complex with gp41 Peptide analog ELD(Orn)WAS
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Deposited on : 2009-07-21
Resolution : 2.24 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
Mogul : 2022.3.0, CSD as543be (2022)
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

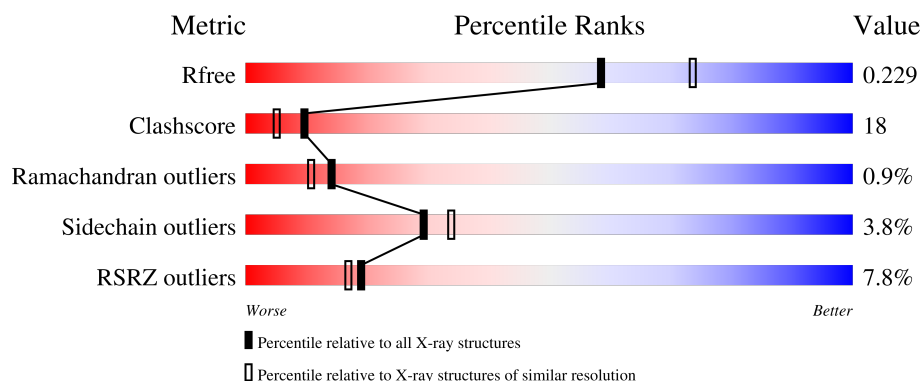
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.24 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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	214	6% 65% 32% .
2	B	237	9% 68% 25% ...
3	C	7	14% 71% 29%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3628 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 2F5 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1644	1025	281	333	5			

- Molecule 2 is a protein called 2F5 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	229	Total	C	N	O	S	0	0	0
			1715	1090	291	327	7			

- Molecule 3 is a protein called gp41 MPER peptide analog.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	7	Total	C	N	O	0	0	0
			59	37	9	13			

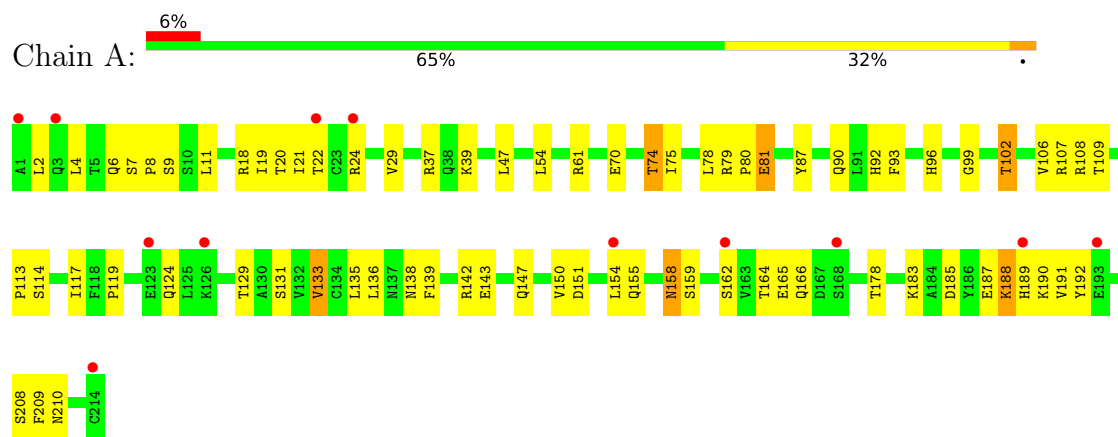
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	99	Total	O	0	0
			99	99		
4	B	102	Total	O	0	0
			102	102		
4	C	9	Total	O	0	0
			9	9		

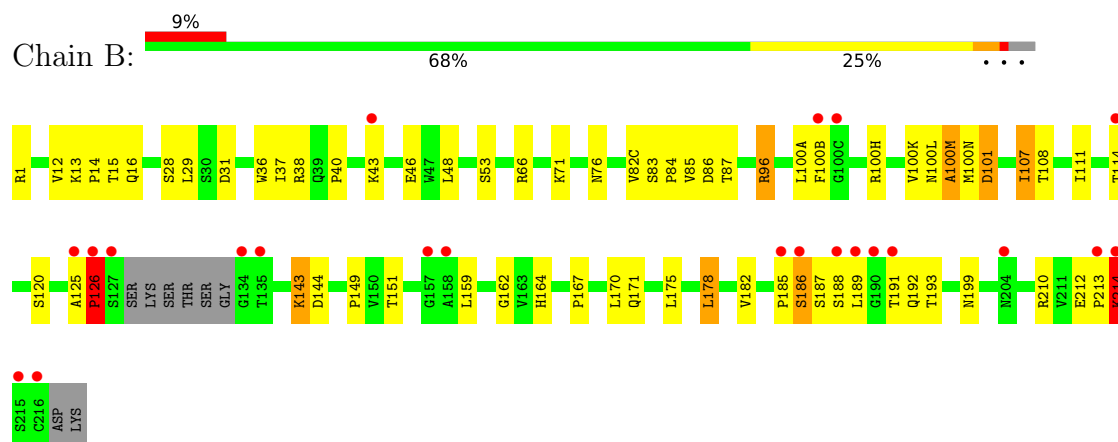
3 Residue-property plots [i](#)

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.

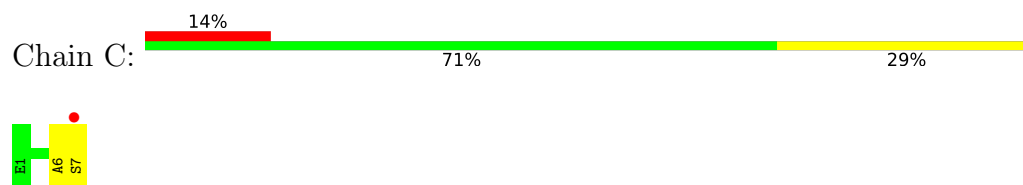
- Molecule 1: 2F5 Fab light chain



- Molecule 2: 2F5 Fab heavy chain



- Molecule 3: gp41 MPER peptide analog



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.40Å 64.90Å 175.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.41 – 2.24 43.41 – 2.24	Depositor EDS
% Data completeness (in resolution range)	92.3 (43.41-2.24) 92.3 (43.41-2.24)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.19 (at 2.22Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.229 , 0.229 0.230 , 0.229	Depositor DCC
R_{free} test set	1543 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3628	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.46% 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

Bond lengths and bond angles in the following residue types are not validated in this section: ORN

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	0.69	0/1681	1.01	5/2284 (0.2%)
2	B	0.70	0/1757	1.13	9/2403 (0.4%)
3	C	0.63	0/51	0.94	0/66
All	All	0.69	0/3489	1.07	14/4753 (0.3%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	125	ALA	CA-C-N	9.04	148.69	127.00
2	B	125	ALA	C-N-CA	9.04	148.69	127.00
2	B	125	ALA	C-N-CD	-8.79	101.26	120.60
2	B	126	PRO	N-CA-C	7.47	131.53	112.10
2	B	143	LYS	N-CA-C	7.06	121.18	109.95
2	B	101	ASP	N-CA-C	6.80	120.79	112.23
2	B	144	ASP	N-CA-C	6.68	120.31	111.28
1	A	114	SER	N-CA-C	-6.11	99.31	109.46
2	B	120	SER	N-CA-C	-6.08	99.28	109.07
1	A	142	ARG	N-CA-C	6.02	117.84	111.28
1	A	96	HIS	N-CA-C	-5.76	102.89	110.43
1	A	133	VAL	N-CA-C	5.61	116.66	108.53
2	B	100(N)	MET	N-CA-C	-5.49	96.80	107.62
1	A	191	VAL	N-CA-C	5.11	115.12	107.51

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	1644	0	1588	65	0
2	B	1715	0	1726	56	0
3	C	59	0	53	2	0
4	A	99	0	0	2	0
4	B	102	0	0	1	0
4	C	9	0	0	0	0
All	All	3628	0	3367	120	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 (120) 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:87:THR:HG22	2:B:111:ILE:H	1.18	1.08
1:A:90:GLN:HE22	1:A:93:PHE:H	1.08	0.93
1:A:20:THR:HG22	1:A:74:THR:HB	1.48	0.92
1:A:190:LYS:HE2	1:A:210:ASN:ND2	1.87	0.88
2:B:86:ASP:HB2	2:B:111:ILE:HD12	1.52	0.87
1:A:90:GLN:NE2	1:A:93:PHE:H	1.73	0.86
1:A:90:GLN:HE21	1:A:92:HIS:H	1.24	0.85
2:B:38:ARG:HD2	2:B:48:LEU:HD21	1.60	0.82
1:A:190:LYS:HE2	1:A:210:ASN:HD22	1.42	0.82
2:B:192:GLN:HE21	2:B:192:GLN:HA	1.45	0.81
1:A:90:GLN:NE2	1:A:92:HIS:H	1.80	0.80
2:B:86:ASP:HB2	2:B:111:ILE:CD1	2.14	0.78
2:B:193:THR:HG22	2:B:210:ARG:HH21	1.46	0.78
1:A:29:VAL:HG11	1:A:90:GLN:HG3	1.65	0.78
2:B:87:THR:HG22	2:B:111:ILE:N	1.99	0.77
2:B:82(C):VAL:HB	2:B:111:ILE:HD13	1.70	0.73
2:B:28:SER:HB3	2:B:31:ASP:OD2	1.89	0.72
2:B:107:ILE:HD13	2:B:108:THR:N	2.03	0.72
1:A:7:SER:HB2	1:A:22:THR:CG2	2.20	0.71
1:A:7:SER:HB2	1:A:22:THR:HG23	1.72	0.70
2:B:188:SER:HA	2:B:191:THR:HG22	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:PRO:O	1:A:102:THR:HB	1.91	0.69
2:B:114:THR:HG21	2:B:175:LEU:HD23	1.74	0.69
2:B:1:ARG:HH11	2:B:1:ARG:HG2	1.59	0.67
2:B:192:GLN:HA	2:B:192:GLN:NE2	2.13	0.64
1:A:106:VAL:H	1:A:166:GLN:HE22	1.42	0.64
2:B:38:ARG:HD3	2:B:46:GLU:OE1	1.98	0.64
1:A:90:GLN:NE2	1:A:92:HIS:N	2.45	0.64
2:B:87:THR:CG2	2:B:111:ILE:H	2.04	0.62
2:B:84:PRO:O	2:B:87:THR:HG23	2.00	0.61
2:B:178:LEU:HD12	2:B:178:LEU:C	2.26	0.60
1:A:29:VAL:HG11	1:A:90:GLN:CG	2.30	0.60
1:A:117:ILE:O	2:B:126:PRO:HG2	2.01	0.60
2:B:191:THR:HG23	2:B:192:GLN:N	2.17	0.60
1:A:79:ARG:HB3	1:A:81:GLU:OE2	2.02	0.59
3:C:6:ALA:O	3:C:7:SER:HB2	2.02	0.59
2:B:40:PRO:HG2	2:B:43:LYS:HB2	1.84	0.59
2:B:100(H):ARG:O	2:B:100(K):VAL:HG22	2.03	0.58
1:A:162:SER:HB3	2:B:167:PRO:HG2	1.86	0.57
1:A:158:ASN:HD22	1:A:158:ASN:H	1.52	0.57
1:A:185:ASP:HA	1:A:188:LYS:NZ	2.19	0.57
2:B:191:THR:HG23	2:B:192:GLN:H	1.70	0.57
1:A:158:ASN:H	1:A:158:ASN:ND2	2.03	0.56
2:B:151:THR:OG1	2:B:199:ASN:HB3	2.06	0.56
1:A:124:GLN:HE22	1:A:131:SER:HB2	1.71	0.56
1:A:183:LYS:O	1:A:187:GLU:HG2	2.06	0.55
2:B:214:LYS:H	2:B:214:LYS:NZ	2.04	0.55
2:B:114:THR:HG21	2:B:175:LEU:CD2	2.37	0.54
1:A:21:ILE:CG2	1:A:102:THR:HG21	2.38	0.54
2:B:36:TRP:C	2:B:37:ILE:HD12	2.33	0.54
1:A:188:LYS:NZ	1:A:188:LYS:HB2	2.23	0.53
2:B:107:ILE:HD13	2:B:107:ILE:C	2.33	0.53
1:A:147:GLN:HG3	1:A:154:LEU:HD13	1.89	0.53
2:B:100(A):LEU:HG	2:B:100(B):PHE:CD1	2.43	0.53
1:A:90:GLN:HE21	1:A:92:HIS:N	1.98	0.53
2:B:29:LEU:O	2:B:53:SER:HB3	2.10	0.52
2:B:29:LEU:O	2:B:53:SER:CB	2.59	0.51
2:B:96:ARG:HD3	2:B:101:ASP:OD2	2.10	0.51
1:A:124:GLN:HG2	1:A:129:THR:O	2.10	0.51
1:A:107:ARG:HG2	1:A:108:ARG:N	2.25	0.51
2:B:143:LYS:NZ	2:B:171:GLN:OE1	2.40	0.50
2:B:193:THR:CG2	2:B:210:ARG:HH21	2.18	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:83:SER:OG	2:B:85:VAL:HG22	2.10	0.50
2:B:213:PRO:HA	2:B:214:LYS:HZ1	1.77	0.50
1:A:113:PRO:HB3	1:A:139:PHE:HB3	1.94	0.50
2:B:37:ILE:HD12	2:B:37:ILE:N	2.27	0.49
1:A:6:GLN:HE21	1:A:99:GLY:HA3	1.77	0.49
2:B:12:VAL:O	2:B:111:ILE:HA	2.13	0.49
1:A:6:GLN:OE1	1:A:102:THR:HG23	2.13	0.49
2:B:100(A):LEU:HG	2:B:100(B):PHE:HD1	1.78	0.49
1:A:155:GLN:HB3	1:A:158:ASN:HD21	1.78	0.48
2:B:38:ARG:HB3	2:B:48:LEU:HD11	1.94	0.48
1:A:21:ILE:HG23	1:A:102:THR:HG21	1.95	0.48
1:A:81:GLU:H	1:A:81:GLU:CD	2.21	0.48
2:B:214:LYS:H	2:B:214:LYS:CE	2.27	0.48
1:A:164:THR:HG22	1:A:165:GLU:N	2.29	0.47
2:B:185:PRO:C	2:B:187:SER:H	2.22	0.47
1:A:158:ASN:ND2	1:A:158:ASN:N	2.58	0.47
1:A:109:THR:HG22	4:A:263:HOH:O	2.14	0.47
1:A:119:PRO:HB3	1:A:209:PHE:CE2	2.50	0.47
3:C:6:ALA:O	3:C:7:SER:CB	2.63	0.47
1:A:159:SER:HA	1:A:178:THR:O	2.14	0.47
1:A:147:GLN:CG	1:A:154:LEU:HD13	2.45	0.46
1:A:7:SER:O	1:A:22:THR:HG22	2.14	0.46
1:A:138:ASN:OD1	2:B:164:HIS:HE1	1.99	0.46
1:A:147:GLN:HG3	1:A:154:LEU:CD1	2.45	0.46
1:A:119:PRO:HB3	1:A:209:PHE:CZ	2.50	0.46
2:B:162:GLY:O	2:B:182:VAL:HA	2.15	0.46
1:A:24:ARG:HG3	1:A:70:GLU:OE2	2.16	0.46
1:A:6:GLN:HE22	1:A:87:TYR:HA	1.80	0.46
2:B:100(L):ASN:O	2:B:100(M):ALA:HB2	2.16	0.46
1:A:18:ARG:HB2	1:A:18:ARG:NH1	2.30	0.46
1:A:19:ILE:HD11	1:A:75:ILE:HD12	1.98	0.45
2:B:29:LEU:HD23	2:B:71:LYS:HB2	1.97	0.45
2:B:76:ASN:ND2	4:B:245:HOH:O	2.45	0.45
1:A:2:LEU:C	1:A:2:LEU:HD23	2.42	0.45
1:A:9:SER:O	1:A:102:THR:HA	2.18	0.44
1:A:143:GLU:CD	1:A:143:GLU:H	2.26	0.44
1:A:79:ARG:HB3	1:A:80:PRO:HD2	2.00	0.44
1:A:108:ARG:HH21	1:A:109:THR:HG23	1.82	0.44
1:A:158:ASN:HD22	1:A:158:ASN:N	2.12	0.44
2:B:210:ARG:NH1	2:B:212:GLU:OE1	2.43	0.43
2:B:192:GLN:NE2	2:B:192:GLN:CA	2.79	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:LEU:C	1:A:78:LEU:HD23	2.43	0.43
2:B:186:SER:HA	2:B:189:LEU:CD1	2.48	0.43
2:B:192:GLN:HE21	2:B:192:GLN:CA	2.17	0.43
1:A:11:LEU:C	1:A:11:LEU:HD23	2.43	0.43
1:A:150:VAL:O	1:A:151:ASP:C	2.62	0.43
2:B:1:ARG:HG2	2:B:1:ARG:NH1	2.32	0.43
1:A:150:VAL:HG11	1:A:189:HIS:ND1	2.33	0.42
2:B:14:PRO:O	2:B:15:THR:OG1	2.35	0.42
1:A:61:ARG:HD3	1:A:79:ARG:NH2	2.34	0.42
1:A:192:TYR:O	1:A:208:SER:HB2	2.19	0.42
1:A:135:LEU:C	1:A:136:LEU:HD12	2.44	0.41
1:A:154:LEU:HD12	1:A:155:GLN:N	2.35	0.41
1:A:39:LYS:HE3	4:A:241:HOH:O	2.21	0.41
2:B:66:ARG:NH2	2:B:86:ASP:OD2	2.51	0.41
2:B:13:LYS:O	2:B:16:GLN:HB2	2.21	0.41
1:A:37:ARG:HB2	1:A:47:LEU:HD11	2.03	0.41
1:A:4:LEU:HD11	1:A:90:GLN:HB2	2.03	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	212/214 (99%)	205 (97%)	7 (3%)	0	100	100
2	B	225/237 (95%)	213 (95%)	8 (4%)	4 (2%)	6	3
3	C	4/7 (57%)	4 (100%)	0	0	100	100
All	All	441/458 (96%)	422 (96%)	15 (3%)	4 (1%)	14	11

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	126	PRO
2	B	214	LYS
2	B	186	SER
2	B	100(M)	ALA

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	187/187 (100%)	180 (96%)	7 (4%)	30	35
2	B	198/207 (96%)	190 (96%)	8 (4%)	28	32
3	C	5/5 (100%)	5 (100%)	0	100	100
All	All	390/399 (98%)	375 (96%)	15 (4%)	29	34

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	LEU
1	A	74	THR
1	A	81	GLU
1	A	102	THR
1	A	133	VAL
1	A	158	ASN
1	A	188	LYS
2	B	96	ARG
2	B	107	ILE
2	B	126	PRO
2	B	149	PRO
2	B	159	LEU
2	B	170	LEU
2	B	178	LEU
2	B	214	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	38	GLN
1	A	90	GLN
1	A	124	GLN
1	A	137	ASN
1	A	147	GLN
1	A	155	GLN
1	A	158	ASN
1	A	166	GLN
1	A	189	HIS
1	A	199	GLN
1	A	210	ASN
2	B	16	GLN
2	B	39	GLN
2	B	76	ASN
2	B	164	HIS
2	B	192	GLN
2	B	199	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ORN	C	4	3	6,7,8	0.71	0	2,7,9	0.44	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ORN	C	4	3	-	0/5/6/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

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	214/214 (100%)	0.40	12 (5%) 30 28	22, 35, 51, 61	0
2	B	229/237 (96%)	0.53	22 (9%) 13 12	21, 34, 59, 83	0
3	C	6/7 (85%)	0.47	1 (16%) 4 3	28, 38, 40, 52	0
All	All	449/458 (98%)	0.47	35 (7%) 19 17	21, 35, 54, 83	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	216	CYS	9.7
2	B	126	PRO	5.6
2	B	215	SER	5.5
2	B	190	GLY	5.3
2	B	127	SER	5.2
2	B	214	LYS	5.0
2	B	100(C)	GLY	4.9
2	B	189	LEU	4.9
2	B	125	ALA	4.7
2	B	114	THR	4.1
2	B	134	GLY	3.9
1	A	154	LEU	3.7
2	B	191	THR	3.5
1	A	1	ALA	3.5
1	A	214	CYS	3.5
2	B	100(B)	PHE	3.4
2	B	135	THR	3.0
2	B	204	ASN	2.9
2	B	188	SER	2.8
1	A	24	ARG	2.8
1	A	193	GLU	2.8
1	A	123	GLU	2.7
1	A	22	THR	2.6

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Mol	Chain	Res	Type	RSRZ
3	C	7	SER	2.5
1	A	126	LYS	2.4
1	A	168	SER	2.4
2	B	43	LYS	2.3
1	A	189	HIS	2.3
2	B	213	PRO	2.3
2	B	158	ALA	2.2
1	A	3	GLN	2.2
2	B	186	SER	2.2
1	A	162	SER	2.1
2	B	157	GLY	2.0
2	B	185	PRO	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ORN	C	4	8/9	0.95	0.08	30,33,36,38	0

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.