



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:28 AM UTC

PDB ID : 1T2K / pdb_00001t2k
Title : Structure Of The DNA Binding Domains Of IRF3, ATF-2 and Jun Bound To DNA
Authors : Panne, D.; Maniatis, T.; Harrison, S.C.
Deposited on : 2004-04-21
Resolution : 3.00 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:

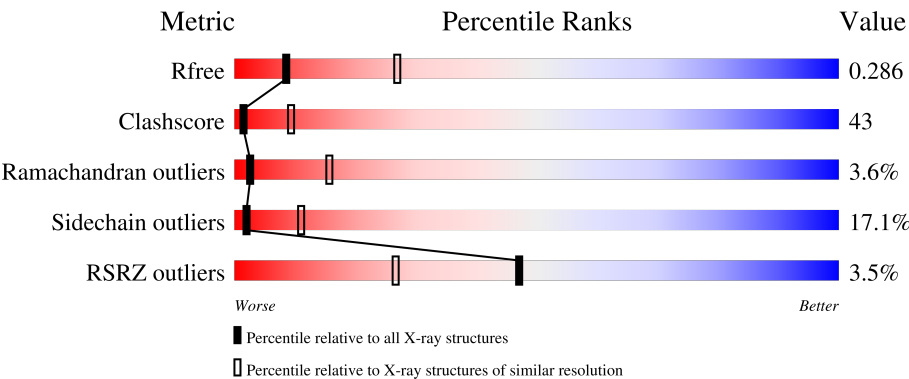
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	E	31	3% 13% 84% .
2	F	31	10% 90%
3	A	112	2% 39% 41% 14% . .
3	B	112	4% 32% 49% 12% . .
4	C	62	10% 53% 34% 13%

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Mol	Chain	Length	Quality of chain
5	D	61	2% 44% 43% 13%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4034 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 DNA chain called 31-MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	31	Total	C	N	O	P	0	0	0
			649	308	139	172	30			

- Molecule 2 is a DNA chain called 31-MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	31	Total	C	N	O	P	0	0	0
			616	301	89	196	30			

- Molecule 3 is a protein called Interferon regulatory factor 3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	A	110	Total	C	N	O	0	0	0
			912	581	172	159			
3	B	108	Total	C	N	O	0	0	0
			888	568	166	154			

- Molecule 4 is a protein called Transcription factor AP-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	C	62	Total	C	N	O	S	0	0	0
			484	293	99	89	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	253	MET	ILE	cloning artifact	UNP P05412
C	269	SER	CYS	engineered mutation	UNP P05412

- Molecule 5 is a protein called Cyclic-AMP-dependent transcription factor ATF-2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	D	61	Total	C	N	O	0	0	0
			477	292	98	87			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	351	SER	CYS	engineered mutation	UNP P15336

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	E	2	Total	O	0	0
			2	2		
6	F	3	Total	O	0	0
			3	3		
6	A	3	Total	O	0	0
			3	3		

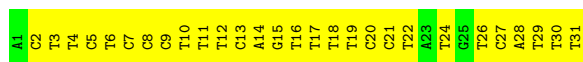
3 Residue-property plots [i](#)

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.

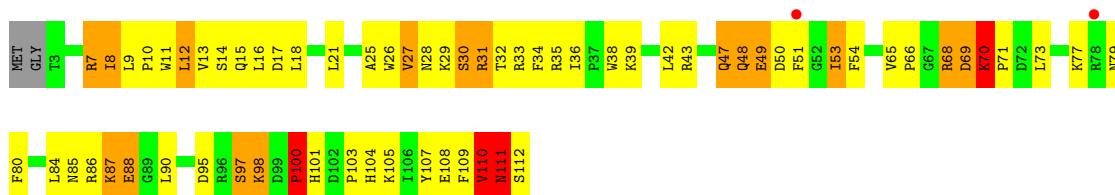
- Molecule 1: 31-MER



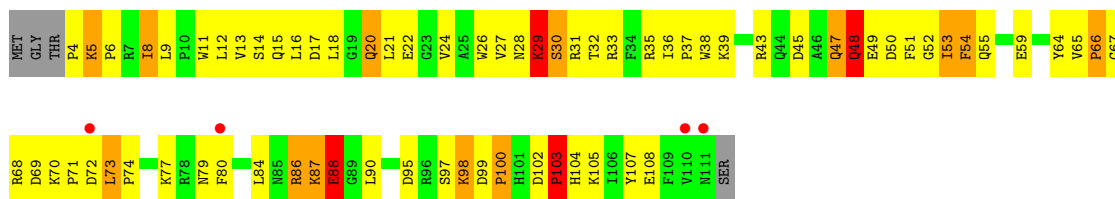
- Molecule 2: 31-MER



- Molecule 3: Interferon regulatory factor 3



- Molecule 3: Interferon regulatory factor 3

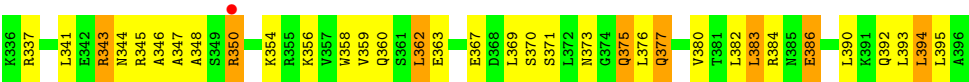


- Molecule 4: Transcription factor AP-1





● Molecule 5: Cyclic-AMP-dependent transcription factor ATF-2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	186.38Å 65.21Å 83.85Å 90.00° 93.44° 90.00°	Depositor
Resolution (Å)	29.98 – 3.00 29.98 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.98-3.00) 98.8 (29.98-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 3.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.257 , 0.296 0.256 , 0.286	Depositor DCC
R_{free} test set	1178 reflections (5.85%)	wwPDB-VP
Wilson B-factor (Å ²)	55.1	Xtriage
Anisotropy	1.286	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 65.3	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	0.94	EDS
Total number of atoms	4034	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.04% 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	E	0.32	0/735	0.65	0/1135
2	F	0.36	0/683	0.82	0/1050
3	A	0.87	2/940 (0.2%)	1.41	18/1273 (1.4%)
3	B	0.73	0/916	1.42	17/1241 (1.4%)
4	C	0.70	0/483	1.24	4/638 (0.6%)
5	D	0.64	0/479	1.22	5/638 (0.8%)
All	All	0.65	2/4236 (0.0%)	1.17	44/5975 (0.7%)

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	E	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	110	VAL	N-CA	5.62	1.53	1.46
3	A	112	SER	C-OXT	-5.45	1.12	1.23

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	103	PRO	CA-N-CD	-13.69	92.83	112.00
3	B	8	ILE	N-CA-C	10.64	121.47	110.62
3	B	48	GLN	OE1-CD-NE2	-9.88	112.72	122.60
3	B	47	GLN	OE1-CD-NE2	-9.82	112.78	122.60
3	A	47	GLN	OE1-CD-NE2	-9.71	112.89	122.60
3	A	48	GLN	OE1-CD-NE2	-9.59	113.01	122.60
3	A	100	PRO	CA-N-CD	-9.01	99.39	112.00
3	A	8	ILE	N-CA-C	8.77	119.56	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	393	LEU	N-CA-C	8.46	120.58	111.36
3	A	36	ILE	N-CA-C	7.59	116.16	107.89
3	A	69	ASP	CA-C-N	7.28	131.39	121.20
3	A	69	ASP	C-N-CA	7.28	131.39	121.20
3	A	101	HIS	N-CA-C	7.15	122.17	113.38
5	D	348	ALA	N-CA-C	-7.12	103.52	111.28
3	B	5	LYS	CA-C-N	7.09	127.08	119.78
3	B	5	LYS	C-N-CA	7.09	127.08	119.78
3	A	47	GLN	CG-CD-NE2	6.69	126.44	116.40
3	B	47	GLN	CG-CD-NE2	6.67	126.40	116.40
3	A	70	LYS	CA-C-N	-6.67	112.79	119.92
3	A	70	LYS	C-N-CA	-6.67	112.79	119.92
3	A	48	GLN	CG-CD-NE2	6.58	126.27	116.40
3	B	48	GLN	N-CA-C	6.55	124.75	110.80
3	B	48	GLN	CG-CD-NE2	6.46	126.09	116.40
3	B	54	PHE	N-CA-C	-6.33	104.38	111.28
3	B	22	GLU	N-CA-C	6.24	119.09	110.35
3	A	27	VAL	N-CA-C	-6.21	106.72	112.43
3	B	103	PRO	CA-C-N	-6.09	111.38	121.80
3	B	103	PRO	C-N-CA	-6.09	111.38	121.80
4	C	261	ARG	N-CA-C	-6.07	104.67	111.28
5	D	394	LEU	N-CA-C	5.95	119.72	112.23
5	D	373	ASN	N-CA-C	-5.62	105.15	111.28
5	D	392	GLN	N-CA-C	5.59	117.18	111.14
3	A	97	SER	N-CA-C	5.56	118.06	111.33
3	B	52	GLY	N-CA-C	5.43	119.72	112.77
3	B	100	PRO	CB-CA-C	5.32	120.34	111.56
3	B	47	GLN	N-CA-C	-5.32	100.99	109.23
4	C	276	ARG	N-CA-C	-5.31	105.49	111.28
3	A	36	ILE	N-CA-CB	-5.22	106.02	111.61
3	B	6	PRO	N-CA-C	5.21	119.49	111.21
3	A	111	ASN	CA-C-N	5.20	131.05	121.70
3	A	111	ASN	C-N-CA	5.20	131.05	121.70
4	C	307	GLN	N-CA-C	5.11	116.93	111.36
4	C	279	ARG	N-CA-C	-5.07	105.75	111.28
3	A	111	ASN	CA-CB-CG	5.02	117.62	112.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	6	DG	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	E	649	0	347	46	0
2	F	616	0	359	70	0
3	A	912	0	893	64	0
3	B	888	0	863	95	0
4	C	484	0	512	31	0
5	D	477	0	483	51	0
6	A	3	0	0	0	0
6	E	2	0	0	0	0
6	F	3	0	0	0	0
All	All	4034	0	3457	323	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 43.

All (323) 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)
5:D:347:ALA:HA	5:D:350:ARG:CG	1.77	1.15
2:F:5:DC:H2''	2:F:6:DT:H5''	1.16	1.08
2:F:29:DT:H2''	2:F:30:DT:H5''	1.33	1.07
3:B:99:ASP:O	3:B:103:PRO:HD3	1.54	1.06
3:A:53:ILE:H	3:A:53:ILE:HD12	1.20	1.04
5:D:347:ALA:HA	5:D:350:ARG:HG3	1.39	1.04
3:B:53:ILE:HD12	3:B:53:ILE:H	1.11	1.04
5:D:347:ALA:HA	5:D:350:ARG:CD	1.89	1.01
2:F:5:DC:C2'	2:F:6:DT:H5''	1.90	1.01
2:F:18:DT:H2''	2:F:19:DT:H5'	1.41	1.00
4:C:280:LEU:HD22	5:D:362:LEU:HD21	1.39	1.00
5:D:347:ALA:HA	5:D:350:ARG:HD2	1.48	0.94
3:B:86:ARG:HB3	3:B:86:ARG:NH1	1.84	0.91
3:B:29:LYS:H	3:B:29:LYS:HD3	1.36	0.89
4:C:280:LEU:CD2	5:D:362:LEU:HD21	2.03	0.89
5:D:347:ALA:CA	5:D:350:ARG:HG3	2.03	0.89
5:D:362:LEU:C	5:D:362:LEU:HD23	1.97	0.88
5:D:346:ALA:O	5:D:350:ARG:HG2	1.73	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:68:ARG:NH1	3:A:69:ASP:OD2	2.08	0.87
3:B:53:ILE:HD12	3:B:53:ILE:N	1.92	0.85
3:B:86:ARG:HB3	3:B:86:ARG:HH11	1.41	0.84
3:B:29:LYS:HD3	3:B:29:LYS:N	1.93	0.84
2:F:5:DC:H2''	2:F:6:DT:C5'	2.03	0.84
3:B:20:GLN:HE21	3:B:20:GLN:HA	1.45	0.81
2:F:29:DT:C2'	2:F:30:DT:H5''	2.10	0.81
2:F:10:DT:H2''	2:F:11:DT:C5'	2.11	0.81
3:B:53:ILE:H	3:B:53:ILE:CD1	1.88	0.81
1:E:3:DA:H2''	1:E:4:DA:H5'	1.61	0.81
2:F:18:DT:H2''	2:F:19:DT:C5'	2.10	0.81
1:E:31:DG:H1	2:F:2:DC:H42	1.27	0.80
2:F:10:DT:H3'	3:B:79:ASN:ND2	1.96	0.80
2:F:18:DT:H6	2:F:18:DT:H5'	1.46	0.80
2:F:29:DT:H2''	2:F:30:DT:C5'	2.11	0.79
1:E:17:DA:H2''	1:E:18:DC:O5'	1.82	0.79
3:A:53:ILE:HD12	3:A:53:ILE:N	1.98	0.78
2:F:10:DT:H2''	2:F:11:DT:H5'	1.66	0.78
4:C:267:SER:O	4:C:271:LYS:HG3	1.84	0.77
5:D:362:LEU:HD23	5:D:363:GLU:N	2.00	0.77
2:F:10:DT:H3'	3:B:79:ASN:HD21	1.46	0.77
3:B:51:PHE:CD2	3:B:73:LEU:HB3	2.20	0.77
4:C:312:VAL:HG21	5:D:394:LEU:HB3	1.65	0.76
3:B:47:GLN:O	3:B:49:GLU:N	2.19	0.75
3:B:9:LEU:O	3:B:13:VAL:HG23	1.86	0.75
2:F:14:DA:H2''	2:F:15:DG:H5''	1.69	0.75
5:D:347:ALA:CA	5:D:350:ARG:CG	2.62	0.75
3:A:9:LEU:O	3:A:13:VAL:HG23	1.86	0.74
5:D:347:ALA:C	5:D:350:ARG:HG3	2.12	0.74
1:E:3:DA:H2''	1:E:4:DA:C5'	2.17	0.74
5:D:347:ALA:O	5:D:350:ARG:HG3	1.88	0.73
3:A:35:ARG:O	3:A:35:ARG:HG3	1.87	0.73
3:A:53:ILE:H	3:A:53:ILE:CD1	1.88	0.72
3:B:28:ASN:CG	3:B:32:THR:HG22	2.14	0.72
1:E:17:DA:H8	1:E:17:DA:H5'	1.56	0.71
3:B:51:PHE:CE2	3:B:73:LEU:HB3	2.26	0.71
2:F:7:DC:H2''	2:F:8:DC:H5'	1.72	0.70
3:B:20:GLN:C	3:B:21:LEU:HD12	2.17	0.70
1:E:7:DA:H2''	1:E:8:DC:H5'	1.74	0.69
5:D:375:GLN:H	5:D:375:GLN:HE21	1.40	0.69
3:A:87:LYS:O	3:A:88:GLU:C	2.36	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:28:DA:H2''	2:F:29:DT:H5''	1.74	0.68
3:B:73:LEU:HB2	3:B:74:PRO:HD3	1.75	0.68
2:F:18:DT:H5'	2:F:18:DT:C6	2.28	0.68
3:A:9:LEU:HD11	3:A:87:LYS:HG3	1.75	0.68
3:A:21:LEU:HD12	3:A:21:LEU:H	1.59	0.68
2:F:3:DT:H2'	2:F:4:DT:H71	1.75	0.67
2:F:12:DT:H4'	2:F:13:DC:OP1	1.93	0.67
4:C:280:LEU:HD13	5:D:362:LEU:HD22	1.76	0.67
2:F:15:DG:H2''	2:F:16:DT:O5'	1.95	0.67
5:D:375:GLN:HE21	5:D:375:GLN:N	1.93	0.66
3:B:43:ARG:HB3	3:B:45:ASP:OD1	1.94	0.66
3:B:73:LEU:HB2	3:B:74:PRO:CD	2.26	0.66
3:B:73:LEU:H	3:B:73:LEU:HD22	1.59	0.66
4:C:312:VAL:CG2	5:D:394:LEU:HB3	2.24	0.66
3:A:21:LEU:HD12	3:A:21:LEU:N	2.10	0.66
2:F:30:DT:H2'	2:F:31:DT:H71	1.77	0.66
3:B:20:GLN:HA	3:B:20:GLN:NE2	2.10	0.65
3:B:35:ARG:HH22	3:B:102:ASP:CG	2.05	0.65
3:B:54:PHE:HD1	3:B:80:PHE:HD2	1.45	0.64
5:D:337:ARG:HG2	5:D:337:ARG:HH11	1.61	0.64
3:B:99:ASP:O	3:B:103:PRO:CD	2.38	0.64
2:F:5:DC:C3'	2:F:6:DT:H5''	2.27	0.63
2:F:16:DT:H3'	3:A:79:ASN:ND2	2.13	0.63
1:E:7:DA:H2''	1:E:8:DC:C5'	2.29	0.63
3:B:38:TRP:CD2	3:B:105:LYS:HD2	2.34	0.62
3:B:69:ASP:O	3:B:70:LYS:C	2.41	0.62
3:A:54:PHE:CD1	3:A:80:PHE:HD2	2.18	0.62
2:F:10:DT:H2''	2:F:11:DT:H5''	1.81	0.62
1:E:28:DG:H1	2:F:5:DC:H42	1.46	0.62
3:B:14:SER:O	3:B:18:LEU:HG	2.00	0.61
3:B:28:ASN:OD1	3:B:28:ASN:O	2.17	0.61
1:E:17:DA:H5'	1:E:17:DA:C8	2.34	0.61
3:B:73:LEU:H	3:B:73:LEU:CD2	2.13	0.61
5:D:371:SER:O	5:D:375:GLN:NE2	2.32	0.61
3:B:54:PHE:HD1	3:B:80:PHE:CD2	2.19	0.61
1:E:7:DA:H1'	1:E:8:DC:H5''	1.83	0.61
3:B:65:VAL:O	3:B:67:GLY:N	2.34	0.61
3:B:73:LEU:HD22	3:B:73:LEU:N	2.14	0.61
4:C:280:LEU:HD13	5:D:362:LEU:CD2	2.31	0.61
5:D:343:ARG:HH11	5:D:343:ARG:HG3	1.66	0.61
2:F:28:DA:C2'	2:F:29:DT:H5''	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:28:ASN:OD1	3:B:28:ASN:C	2.44	0.60
3:A:110:VAL:O	3:A:111:ASN:HB3	2.01	0.60
3:B:80:PHE:O	3:B:80:PHE:HD1	1.85	0.60
3:A:98:LYS:O	3:A:100:PRO:HD3	2.01	0.60
2:F:6:DT:H2''	2:F:7:DC:C6	2.36	0.60
3:A:39:LYS:HE2	3:A:50:ASP:CG	2.26	0.60
2:F:27:DC:OP2	5:D:337:ARG:NH2	2.34	0.59
3:A:47:GLN:HG3	3:A:49:GLU:HB3	1.85	0.59
4:C:287:LEU:O	4:C:291:ASN:HB3	2.03	0.59
3:A:73:LEU:N	3:A:73:LEU:HD22	2.17	0.58
4:C:261:ARG:HG2	4:C:261:ARG:HH11	1.68	0.58
1:E:8:DC:H6	4:C:270:ARG:NH1	2.02	0.58
2:F:28:DA:H2''	2:F:29:DT:C5'	2.32	0.58
5:D:377:GLN:HE21	5:D:377:GLN:HA	1.67	0.58
2:F:26:DT:H2''	2:F:27:DC:O5'	2.04	0.58
2:F:20:DC:H2''	2:F:21:DC:C5'	2.34	0.57
5:D:362:LEU:C	5:D:362:LEU:CD2	2.69	0.57
1:E:23:DA:H2''	1:E:24:DG:H5''	1.86	0.57
1:E:26:DG:H2''	1:E:27:DA:C8	2.39	0.57
1:E:14:DA:H2''	1:E:15:DA:H5'	1.86	0.57
5:D:356:LYS:O	5:D:360:GLN:HG3	2.05	0.57
3:A:48:GLN:NE2	3:A:49:GLU:N	2.53	0.56
5:D:347:ALA:CA	5:D:350:ARG:HD2	2.28	0.56
3:A:48:GLN:NE2	3:A:49:GLU:H	2.03	0.56
3:A:38:TRP:NE1	3:A:77:LYS:HE3	2.19	0.56
2:F:20:DC:H1'	2:F:21:DC:H5''	1.88	0.56
4:C:291:ASN:HD22	4:C:292:SER:N	2.04	0.56
4:C:299:ASN:O	4:C:303:GLU:HG3	2.06	0.56
2:F:30:DT:H2''	2:F:31:DT:O5'	2.06	0.55
3:A:73:LEU:HD22	3:A:73:LEU:H	1.70	0.55
3:B:35:ARG:NH2	3:B:102:ASP:OD2	2.33	0.55
4:C:274:LEU:HD23	4:C:277:ILE:HD12	1.88	0.55
1:E:19:DT:H2''	1:E:20:DG:C8	2.41	0.55
3:B:80:PHE:CD1	3:B:80:PHE:C	2.83	0.55
4:C:273:LYS:O	4:C:277:ILE:HG13	2.06	0.55
5:D:377:GLN:HE21	5:D:377:GLN:CA	2.20	0.55
3:A:110:VAL:O	3:A:111:ASN:CB	2.54	0.55
3:B:16:LEU:HB3	3:B:26:TRP:NE1	2.22	0.55
1:E:5:DT:OP1	5:D:354:LYS:NZ	2.36	0.55
3:A:38:TRP:NE1	3:A:77:LYS:CE	2.70	0.55
4:C:280:LEU:HD11	5:D:363:GLU:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:28:DG:H2''	1:E:29:DA:C8	2.43	0.54
1:E:9:DA:H1'	1:E:10:DT:H5''	1.89	0.54
3:B:29:LYS:H	3:B:29:LYS:CD	2.16	0.54
4:C:304:GLN:HA	4:C:307:GLN:HB3	1.90	0.54
3:B:29:LYS:C	3:B:31:ARG:H	2.15	0.54
3:A:14:SER:O	3:A:18:LEU:HG	2.08	0.53
3:B:68:ARG:O	3:B:68:ARG:HD3	2.07	0.53
3:B:33:ARG:HG2	3:B:108:GLU:HB2	1.90	0.53
3:A:54:PHE:CD1	3:A:80:PHE:CD2	2.96	0.53
1:E:29:DA:H2''	1:E:30:DA:H5'	1.90	0.53
3:B:65:VAL:O	3:B:69:ASP:HB2	2.09	0.53
4:C:287:LEU:HD23	5:D:369:LEU:CB	2.39	0.53
1:E:25:DG:H2'	1:E:26:DG:C8	2.44	0.53
2:F:3:DT:H5'	2:F:3:DT:H6	1.74	0.53
3:B:64:TYR:O	3:B:66:PRO:HD3	2.08	0.52
3:B:98:LYS:O	3:B:100:PRO:HD3	2.09	0.52
3:A:9:LEU:HD11	3:A:87:LYS:CG	2.39	0.52
2:F:20:DC:H2''	2:F:21:DC:H5''	1.92	0.52
3:B:33:ARG:NH1	3:B:108:GLU:OE1	2.43	0.52
3:A:54:PHE:HD1	3:A:80:PHE:CD2	2.27	0.52
4:C:291:ASN:C	4:C:291:ASN:ND2	2.67	0.52
3:A:7:ARG:O	3:A:10:PRO:HD2	2.09	0.52
1:E:8:DC:C6	4:C:270:ARG:NH1	2.77	0.52
2:F:20:DC:C2'	2:F:21:DC:H5''	2.40	0.52
5:D:346:ALA:O	5:D:350:ARG:CG	2.52	0.52
2:F:10:DT:C6	2:F:11:DT:H72	2.46	0.51
2:F:27:DC:C2'	2:F:28:DA:C8	2.93	0.51
1:E:4:DA:H2''	1:E:5:DT:O5'	2.11	0.51
3:B:38:TRP:NE1	3:B:77:LYS:HE3	2.24	0.51
3:B:27:VAL:HG12	3:B:27:VAL:O	2.10	0.51
3:B:68:ARG:HA	3:B:68:ARG:NE	2.25	0.51
3:A:73:LEU:H	3:A:73:LEU:CD2	2.23	0.51
5:D:390:LEU:HD22	5:D:394:LEU:HD13	1.91	0.51
2:F:18:DT:OP1	3:B:98:LYS:HE2	2.10	0.51
5:D:390:LEU:HD22	5:D:394:LEU:CD1	2.41	0.51
1:E:18:DC:H2''	1:E:19:DT:H5'	1.93	0.51
3:A:54:PHE:CE1	3:A:80:PHE:HD2	2.29	0.51
1:E:2:DA:H2''	1:E:3:DA:OP2	2.11	0.50
3:B:32:THR:HG23	3:B:33:ARG:HG3	1.93	0.50
2:F:27:DC:H2'	2:F:28:DA:C8	2.47	0.50
3:B:84:LEU:HD22	3:B:90:LEU:CD2	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:10:DT:H5''	3:B:79:ASN:HD22	1.76	0.50
5:D:377:GLN:HA	5:D:377:GLN:NE2	2.27	0.50
3:B:33:ARG:HG2	3:B:108:GLU:HG3	1.93	0.49
3:B:33:ARG:CG	3:B:108:GLU:HG3	2.41	0.49
3:B:28:ASN:O	3:B:29:LYS:C	2.55	0.49
3:A:51:PHE:CE2	3:A:73:LEU:HB3	2.48	0.49
3:B:48:GLN:NE2	3:B:55:GLN:HE22	2.10	0.49
3:B:80:PHE:HD1	3:B:80:PHE:C	2.20	0.49
2:F:18:DT:H2'	2:F:19:DT:H71	1.94	0.49
3:B:36:ILE:O	3:B:104:HIS:HB2	2.13	0.49
3:B:86:ARG:HB3	3:B:86:ARG:CZ	2.42	0.48
4:C:267:SER:OG	4:C:271:LYS:HE3	2.12	0.48
1:E:25:DG:H4'	1:E:25:DG:OP1	2.13	0.48
3:A:84:LEU:HD22	3:A:90:LEU:CD2	2.43	0.48
1:E:24:DG:H2''	1:E:25:DG:O5'	2.14	0.48
3:B:38:TRP:CE3	3:B:105:LYS:HD2	2.48	0.48
4:C:311:LYS:HB3	5:D:395:LEU:HD11	1.94	0.48
2:F:9:DC:OP2	3:B:87:LYS:HE2	2.12	0.48
3:A:87:LYS:O	3:A:88:GLU:O	2.31	0.48
3:B:73:LEU:CB	3:B:74:PRO:HD3	2.44	0.48
4:C:257:ARG:O	4:C:257:ARG:HG3	2.14	0.48
1:E:31:DG:H1	2:F:2:DC:N4	2.02	0.48
3:A:25:ALA:O	3:A:35:ARG:HG2	2.14	0.47
1:E:8:DC:H2'	4:C:270:ARG:HH12	1.80	0.47
5:D:359:VAL:O	5:D:363:GLU:HG3	2.13	0.47
2:F:10:DT:C3'	3:B:79:ASN:ND2	2.72	0.47
4:C:255:ALA:O	4:C:259:ARG:HG2	2.13	0.47
1:E:13:DG:H2''	1:E:14:DA:C8	2.49	0.47
1:E:28:DG:H1	2:F:5:DC:N4	2.12	0.47
2:F:9:DC:H2'	2:F:10:DT:H72	1.97	0.47
2:F:14:DA:C2'	2:F:15:DG:H5''	2.41	0.47
2:F:27:DC:H2''	2:F:28:DA:C8	2.49	0.47
3:A:21:LEU:H	3:A:21:LEU:CD1	2.26	0.47
4:C:291:ASN:ND2	4:C:292:SER:N	2.62	0.47
4:C:280:LEU:CD1	5:D:362:LEU:CD2	2.93	0.47
2:F:27:DC:H5	5:D:344:ASN:OD1	1.98	0.47
1:E:26:DG:H2''	1:E:27:DA:N7	2.30	0.47
3:B:11:TRP:O	3:B:15:GLN:HG2	2.15	0.47
3:B:21:LEU:HD12	3:B:21:LEU:N	2.30	0.47
3:B:33:ARG:HG2	3:B:108:GLU:CB	2.44	0.46
3:A:11:TRP:O	3:A:15:GLN:HG2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:24:DT:OP1	3:A:98:LYS:HE2	2.15	0.46
3:B:31:ARG:NH1	3:B:31:ARG:HG3	2.31	0.46
3:B:73:LEU:CD2	3:B:73:LEU:N	2.77	0.46
4:C:260:MET:HG2	4:C:263:ARG:NH2	2.31	0.46
4:C:287:LEU:HD23	5:D:369:LEU:HB2	1.98	0.46
3:A:39:LYS:HD2	3:A:43:ARG:HD3	1.97	0.45
2:F:15:DG:OP1	3:A:7:ARG:HA	2.17	0.45
3:B:73:LEU:CB	3:B:74:PRO:CD	2.90	0.45
2:F:14:DA:H1'	2:F:15:DG:O4'	2.15	0.45
3:B:24:VAL:HG21	3:B:53:ILE:CG1	2.45	0.45
1:E:12:DG:H2''	1:E:13:DG:OP2	2.17	0.45
1:E:6:DG:H2''	1:E:7:DA:H5'	1.99	0.45
1:E:17:DA:H8	1:E:17:DA:C5'	2.27	0.45
3:B:31:ARG:HG3	3:B:31:ARG:HH11	1.82	0.45
3:B:80:PHE:HE1	3:B:84:LEU:HG	1.80	0.45
2:F:2:DC:H2''	2:F:3:DT:H5''	1.98	0.45
2:F:10:DT:C3'	3:B:79:ASN:HD21	2.23	0.45
2:F:18:DT:C2'	2:F:19:DT:C5'	2.89	0.45
3:B:30:SER:O	3:B:31:ARG:C	2.58	0.45
2:F:7:DC:H1'	2:F:8:DC:H5''	1.99	0.45
3:A:87:LYS:HA	3:A:87:LYS:HD2	1.72	0.45
3:A:33:ARG:HG3	3:A:108:GLU:HG3	1.99	0.45
3:B:99:ASP:O	3:B:100:PRO:C	2.59	0.45
1:E:17:DA:C8	1:E:17:DA:C5'	3.00	0.45
2:F:24:DT:H6	2:F:24:DT:H5'	1.82	0.45
3:A:98:LYS:HE3	3:A:98:LYS:HB3	1.71	0.45
3:B:33:ARG:HG2	3:B:108:GLU:CG	2.47	0.45
5:D:380:VAL:O	5:D:384:ARG:HG3	2.16	0.45
1:E:12:DG:H2''	1:E:13:DG:H5'	1.99	0.44
3:A:109:PHE:C	3:A:110:VAL:HG13	2.41	0.44
3:B:39:LYS:HA	3:B:103:PRO:CB	2.47	0.44
4:C:287:LEU:HD23	5:D:370:SER:N	2.32	0.44
2:F:16:DT:C6	2:F:17:DT:H72	2.53	0.44
2:F:21:DC:H2''	2:F:22:DT:H6	1.82	0.44
1:E:12:DG:H1'	1:E:13:DG:H5''	2.00	0.44
1:E:23:DA:H2''	1:E:24:DG:C5'	2.48	0.44
3:A:103:PRO:O	3:A:104:HIS:HB3	2.18	0.44
3:B:87:LYS:O	3:B:88:GLU:C	2.61	0.44
3:A:28:ASN:CG	3:A:32:THR:HG22	2.43	0.43
3:B:59:GLU:HG2	3:B:64:TYR:O	2.19	0.43
1:E:7:DA:C1'	1:E:8:DC:H5''	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:10:DT:H2''	1:E:11:DA:H8	1.84	0.43
2:F:16:DT:H3'	3:A:79:ASN:HD21	1.84	0.43
5:D:390:LEU:HD13	5:D:390:LEU:O	2.18	0.43
1:E:9:DA:C2'	1:E:10:DT:H5''	2.49	0.43
3:B:72:ASP:O	3:B:73:LEU:C	2.61	0.43
3:B:16:LEU:HB3	3:B:26:TRP:CE2	2.53	0.43
5:D:343:ARG:HH11	5:D:343:ARG:CG	2.31	0.43
1:E:3:DA:H2''	1:E:4:DA:H5''	1.96	0.43
3:A:12:LEU:O	3:A:12:LEU:HD22	2.18	0.43
3:A:21:LEU:N	3:A:21:LEU:CD1	2.80	0.42
3:A:42:LEU:HD23	3:A:42:LEU:HA	1.85	0.42
3:B:90:LEU:HD11	3:B:107:TYR:HB3	2.01	0.42
3:A:38:TRP:CD2	3:A:105:LYS:HD2	2.53	0.42
5:D:383:LEU:HA	5:D:386:GLU:HB3	2.01	0.42
3:A:27:VAL:HG12	3:A:27:VAL:O	2.19	0.42
3:A:38:TRP:CE3	3:A:105:LYS:HD2	2.54	0.42
3:B:98:LYS:HB3	3:B:98:LYS:HE3	1.61	0.42
5:D:390:LEU:HD13	5:D:390:LEU:C	2.44	0.42
2:F:27:DC:H4'	2:F:27:DC:OP1	2.20	0.42
3:A:70:LYS:H	3:A:70:LYS:HG2	1.34	0.42
5:D:390:LEU:O	5:D:394:LEU:HB2	2.18	0.42
2:F:29:DT:H2'	2:F:30:DT:H71	2.01	0.42
3:B:51:PHE:CD2	3:B:73:LEU:HD12	2.54	0.42
1:E:7:DA:C2'	1:E:8:DC:C5'	2.97	0.42
2:F:5:DC:H4'	2:F:5:DC:OP1	2.19	0.42
3:A:68:ARG:HG2	3:A:69:ASP:CG	2.45	0.42
3:B:65:VAL:C	3:B:67:GLY:N	2.77	0.42
4:C:280:LEU:O	4:C:280:LEU:HD23	2.19	0.42
1:E:19:DT:H2''	1:E:20:DG:N7	2.35	0.42
2:F:21:DC:H2''	2:F:22:DT:C6	2.54	0.42
5:D:337:ARG:NH2	5:D:341:LEU:HD11	2.35	0.42
3:B:95:ASP:OD2	3:B:97:SER:HB2	2.19	0.41
2:F:13:DC:H2''	2:F:14:DA:O5'	2.19	0.41
3:A:34:PHE:CE1	3:A:107:TYR:HB2	2.55	0.41
3:B:26:TRP:CG	3:B:31:ARG:HD2	2.55	0.41
1:E:14:DA:H1'	1:E:15:DA:H5''	2.03	0.41
3:A:95:ASP:OD2	3:A:97:SER:CB	2.69	0.41
1:E:14:DA:H2''	1:E:15:DA:C5'	2.48	0.41
3:A:16:LEU:O	3:A:17:ASP:C	2.62	0.41
3:B:80:PHE:CE1	3:B:84:LEU:HG	2.55	0.41
5:D:337:ARG:HG2	5:D:337:ARG:NH1	2.31	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:16:DT:H5''	2:F:16:DT:H6	1.84	0.41
2:F:15:DG:H2'	2:F:16:DT:H72	2.03	0.41
3:A:16:LEU:HB3	3:A:26:TRP:NE1	2.36	0.41
3:B:28:ASN:O	3:B:31:ARG:N	2.54	0.41
3:B:36:ILE:HA	3:B:37:PRO:HD3	1.91	0.41
3:B:65:VAL:O	3:B:66:PRO:C	2.64	0.41
3:A:111:ASN:O	3:A:111:ASN:ND2	2.54	0.41
5:D:362:LEU:HD23	5:D:363:GLU:CA	2.50	0.41
3:A:65:VAL:O	3:A:66:PRO:C	2.64	0.40
3:A:70:LYS:O	3:A:71:PRO:C	2.60	0.40
3:B:4:PRO:O	3:B:5:LYS:HB3	2.21	0.40
2:F:20:DC:OP1	3:B:43:ARG:HA	2.21	0.40
3:A:26:TRP:CE3	3:A:31:ARG:HA	2.56	0.40
4:C:261:ARG:HG2	4:C:261:ARG:NH1	2.34	0.40
3:A:30:SER:O	3:A:31:ARG:C	2.65	0.40
3:A:48:GLN:O	3:A:49:GLU:C	2.65	0.40
3:B:17:ASP:OD1	3:B:31:ARG:NH1	2.51	0.40
5:D:376:LEU:O	5:D:380:VAL:HG23	2.22	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	
3	A	108/112 (96%)	89 (82%)	13 (12%)	6 (6%)	1	8
3	B	106/112 (95%)	88 (83%)	12 (11%)	6 (6%)	1	8
4	C	60/62 (97%)	59 (98%)	1 (2%)	0	100	100
5	D	59/61 (97%)	58 (98%)	1 (2%)	0	100	100
All	All	333/347 (96%)	294 (88%)	27 (8%)	12 (4%)	2	16

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	30	SER
3	A	88	GLU
3	B	48	GLN
3	A	111	ASN
3	B	29	LYS
3	B	66	PRO
3	A	31	ARG
3	B	88	GLU
3	A	29	LYS
3	B	30	SER
3	B	71	PRO
3	A	110	VAL

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	
3	A	94/96 (98%)	81 (86%)	13 (14%)	3	17
3	B	90/96 (94%)	77 (86%)	13 (14%)	3	15
4	C	49/54 (91%)	38 (78%)	11 (22%)	1	5
5	D	47/54 (87%)	36 (77%)	11 (23%)	1	4
All	All	280/300 (93%)	232 (83%)	48 (17%)	2	11

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	7	ARG
3	A	8	ILE
3	A	12	LEU
3	A	49	GLU
3	A	53	ILE
3	A	68	ARG
3	A	70	LYS
3	A	85	ASN
3	A	86	ARG
3	A	87	LYS

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Mol	Chain	Res	Type
3	A	98	LYS
3	A	100	PRO
3	A	111	ASN
3	B	8	ILE
3	B	12	LEU
3	B	20	GLN
3	B	29	LYS
3	B	48	GLN
3	B	50	ASP
3	B	53	ILE
3	B	73	LEU
3	B	86	ARG
3	B	87	LYS
3	B	88	GLU
3	B	98	LYS
3	B	103	PRO
4	C	256	GLU
4	C	257	ARG
4	C	259	ARG
4	C	262	ASN
4	C	277	ILE
4	C	280	LEU
4	C	291	ASN
4	C	296	SER
4	C	299	ASN
4	C	300	MET
4	C	305	VAL
5	D	343	ARG
5	D	345	ARG
5	D	350	ARG
5	D	358	TRP
5	D	362	LEU
5	D	367	GLU
5	D	375	GLN
5	D	377	GLN
5	D	382	LEU
5	D	383	LEU
5	D	386	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	40	HIS
3	A	47	GLN
3	A	48	GLN
3	A	111	ASN
3	B	20	GLN
3	B	44	GLN
3	B	48	GLN
3	B	55	GLN
3	B	79	ASN
4	C	262	ASN
4	C	291	ASN
4	C	314	ASN
5	D	375	GLN
5	D	377	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 ⓘ

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	E	31/31 (100%)	0.30	1 (3%) 50 29	49, 74, 95, 98	0
2	F	31/31 (100%)	0.25	0 100 100	50, 74, 103, 117	0
3	A	110/112 (98%)	0.34	2 (1%) 67 44	44, 66, 97, 102	0
3	B	108/112 (96%)	0.75	4 (3%) 45 25	67, 101, 118, 128	0
4	C	62/62 (100%)	0.53	6 (9%) 13 7	96, 122, 137, 145	0
5	D	61/61 (100%)	0.31	1 (1%) 70 47	47, 109, 133, 145	0
All	All	403/409 (98%)	0.47	14 (3%) 47 27	44, 93, 131, 145	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	110	VAL	3.3
4	C	297	THR	2.9
5	D	350	ARG	2.5
3	B	72	ASP	2.5
3	B	111	ASN	2.4
4	C	309	LYS	2.3
4	C	298	ALA	2.3
4	C	313	MET	2.3
4	C	308	LEU	2.1
1	E	1	DT	2.1
4	C	301	LEU	2.1
3	A	78	ARG	2.0
3	A	51	PHE	2.0
3	B	80	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 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.