



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:05 AM UTC

PDB ID : 8EJR / pdb_00008ejr
Title : Kelch domain of human KEAP1 bound to Nrf2 linear peptide, Ac-GDPETGE-NH2
Authors : Muellers, S.N.; Allen, K.N.
Deposited on : 2022-09-18
Resolution : 2.08 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

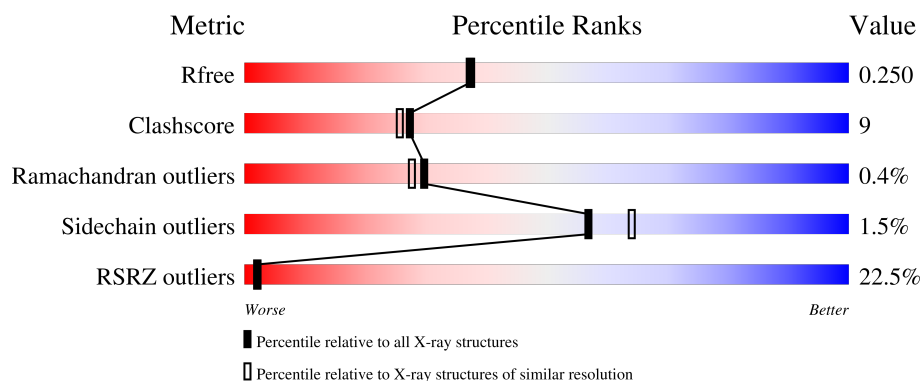
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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	336	6% 75% 9% 16%
1	B	336	32% 62% 20% 16%
2	C	9	22% 89% 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4674 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 Kelch-like ECH-associated protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	S	0	0	0
			2170	1351	392	412	15			
1	B	283	Total	C	N	O	S	0	0	0
			2170	1351	392	412	15			

There are 58 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	289	MET	-	expression tag	UNP Q14145
A	290	GLY	-	expression tag	UNP Q14145
A	291	SER	-	expression tag	UNP Q14145
A	292	SER	-	expression tag	UNP Q14145
A	293	HIS	-	expression tag	UNP Q14145
A	294	HIS	-	expression tag	UNP Q14145
A	295	HIS	-	expression tag	UNP Q14145
A	296	HIS	-	expression tag	UNP Q14145
A	297	HIS	-	expression tag	UNP Q14145
A	298	HIS	-	expression tag	UNP Q14145
A	299	SER	-	expression tag	UNP Q14145
A	300	SER	-	expression tag	UNP Q14145
A	301	GLY	-	expression tag	UNP Q14145
A	302	GLY	-	expression tag	UNP Q14145
A	303	GLU	-	expression tag	UNP Q14145
A	304	ASN	-	expression tag	UNP Q14145
A	305	LEU	-	expression tag	UNP Q14145
A	306	TYR	-	expression tag	UNP Q14145
A	307	PHE	-	expression tag	UNP Q14145
A	308	GLN	-	expression tag	UNP Q14145
A	309	GLY	-	expression tag	UNP Q14145
A	310	HIS	-	expression tag	UNP Q14145
A	311	MET	-	expression tag	UNP Q14145
A	319	SER	CYS	conflict	UNP Q14145
A	540	ALA	GLU	engineered mutation	UNP Q14145

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Chain	Residue	Modelled	Actual	Comment	Reference
A	542	ALA	GLU	engineered mutation	UNP Q14145
A	613	SER	CYS	engineered mutation	UNP Q14145
A	622	SER	CYS	conflict	UNP Q14145
A	624	SER	CYS	conflict	UNP Q14145
B	289	MET	-	expression tag	UNP Q14145
B	290	GLY	-	expression tag	UNP Q14145
B	291	SER	-	expression tag	UNP Q14145
B	292	SER	-	expression tag	UNP Q14145
B	293	HIS	-	expression tag	UNP Q14145
B	294	HIS	-	expression tag	UNP Q14145
B	295	HIS	-	expression tag	UNP Q14145
B	296	HIS	-	expression tag	UNP Q14145
B	297	HIS	-	expression tag	UNP Q14145
B	298	HIS	-	expression tag	UNP Q14145
B	299	SER	-	expression tag	UNP Q14145
B	300	SER	-	expression tag	UNP Q14145
B	301	GLY	-	expression tag	UNP Q14145
B	302	GLY	-	expression tag	UNP Q14145
B	303	GLU	-	expression tag	UNP Q14145
B	304	ASN	-	expression tag	UNP Q14145
B	305	LEU	-	expression tag	UNP Q14145
B	306	TYR	-	expression tag	UNP Q14145
B	307	PHE	-	expression tag	UNP Q14145
B	308	GLN	-	expression tag	UNP Q14145
B	309	GLY	-	expression tag	UNP Q14145
B	310	HIS	-	expression tag	UNP Q14145
B	311	MET	-	expression tag	UNP Q14145
B	319	SER	CYS	conflict	UNP Q14145
B	540	ALA	GLU	engineered mutation	UNP Q14145
B	542	ALA	GLU	engineered mutation	UNP Q14145
B	613	SER	CYS	engineered mutation	UNP Q14145
B	622	SER	CYS	conflict	UNP Q14145
B	624	SER	CYS	conflict	UNP Q14145

- Molecule 2 is a protein called Linear peptide from Nuclear factor erythroid 2-related factor 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	9	Total	C	N	O	0	0	1
			52	29	8	15			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	75	ACE	-	acetylation	UNP Q16236
C	76	GLY	LEU	engineered mutation	UNP Q16236
C	78	PRO	GLU	engineered mutation	UNP Q16236
C	83	NH2	-	amidation	UNP Q16236

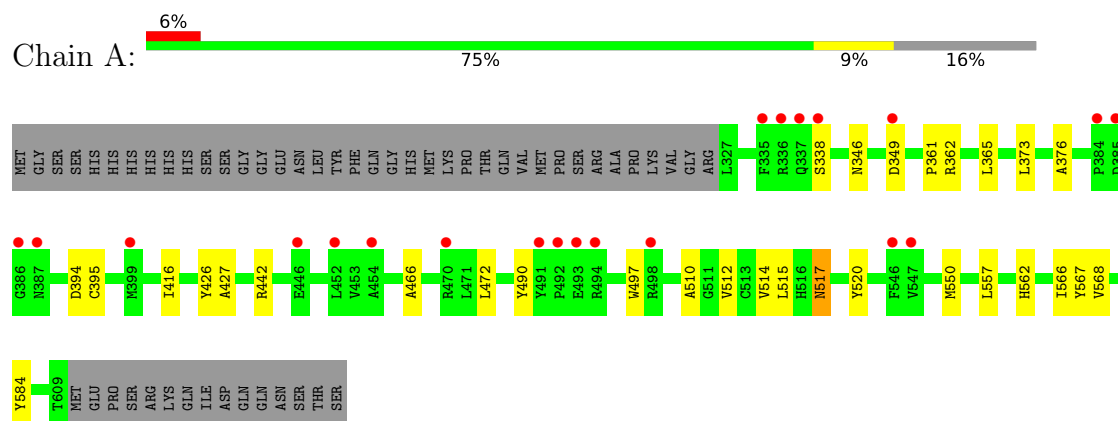
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	179	Total	O	0	0
			179	179		
3	B	96	Total	O	0	0
			96	96		
3	C	7	Total	O	0	0
			7	7		

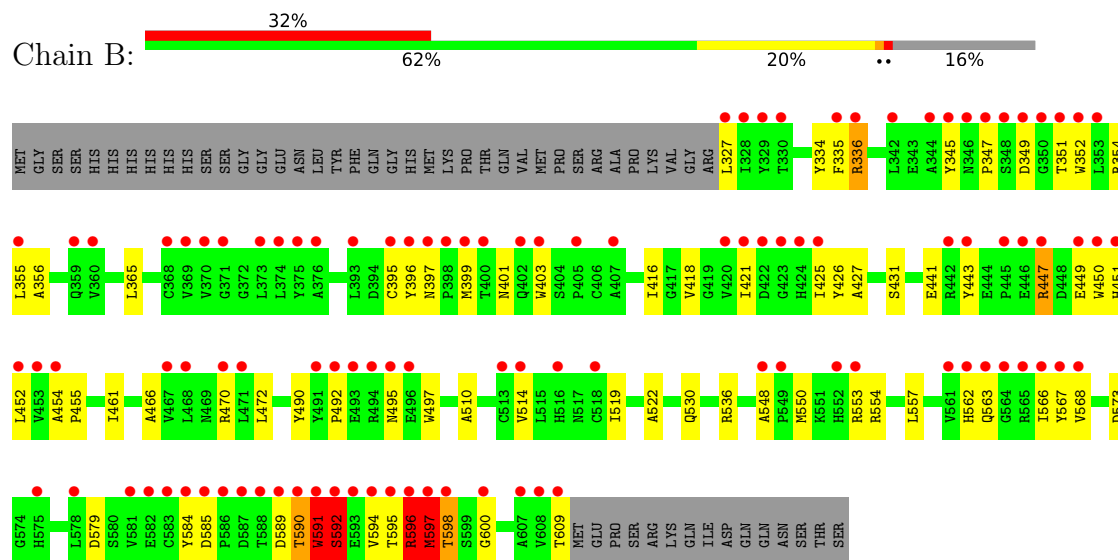
3 Residue-property plots

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.

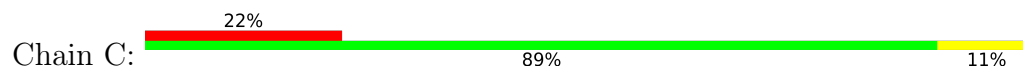
- Molecule 1: Kelch-like ECH-associated protein 1

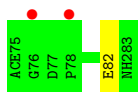


- Molecule 1: Kelch-like ECH-associated protein 1



- Molecule 2: Linear peptide from Nuclear factor erythroid 2-related factor 2





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	162.81Å 68.91Å 77.59Å 90.00° 117.66° 90.00°	Depositor
Resolution (Å)	29.41 – 2.08 29.41 – 2.08	Depositor EDS
% Data completeness (in resolution range)	99.0 (29.41-2.08) 98.9 (29.41-2.08)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 2.08Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.222 , 0.249 0.223 , 0.250	Depositor DCC
R_{free} test set	1993 reflections (4.14%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.567	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.010 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4674	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.39% 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: ACE, NH2

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	0.40	1/2223 (0.0%)	0.57	0/3029
1	B	0.79	9/2223 (0.4%)	1.00	14/3029 (0.5%)
2	C	0.13	0/49	0.31	0/66
All	All	0.63	10/4495 (0.2%)	0.81	14/6124 (0.2%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	591	TRP	CA-C	14.18	1.71	1.52
1	B	592	SER	N-CA	11.02	1.60	1.46
1	B	592	SER	CA-C	10.25	1.65	1.52
1	B	510	ALA	C-N	9.52	1.38	1.33
1	B	591	TRP	N-CA	8.31	1.57	1.46
1	B	590	THR	CA-C	6.83	1.61	1.52
1	A	510	ALA	C-N	5.45	1.36	1.33
1	B	591	TRP	C-N	5.45	1.40	1.33
1	B	590	THR	C-N	5.20	1.40	1.33
1	B	590	THR	N-CA	5.12	1.52	1.46

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	591	TRP	N-CA-C	14.63	129.49	108.86
1	B	592	SER	N-CA-C	13.01	126.70	108.74
1	B	585	ASP	CA-C-N	10.76	130.16	118.97
1	B	585	ASP	C-N-CA	10.76	130.16	118.97
1	B	447	ARG	CG-CD-NE	-9.81	90.42	112.00
1	B	597	MET	CA-C-N	9.47	139.62	121.54
1	B	597	MET	C-N-CA	9.47	139.62	121.54
1	B	590	THR	N-CA-C	7.75	119.44	108.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	597	MET	N-CA-C	6.50	124.65	110.80
1	B	596	ARG	N-CA-C	-6.36	102.90	110.41
1	B	590	THR	CA-C-O	-6.07	114.93	121.36
1	B	492	PRO	CA-C-N	5.80	129.13	120.31
1	B	492	PRO	C-N-CA	5.80	129.13	120.31
1	B	598	THR	N-CA-CB	5.27	119.39	110.49

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	2170	0	2066	20	1
1	B	2170	0	2066	58	1
2	C	52	0	39	1	0
3	A	179	0	0	0	0
3	B	96	0	0	5	0
3	C	7	0	0	0	0
All	All	4674	0	4171	78	1

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 9.

All (78) 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:A:550:MET:HE3	1:A:568:VAL:HG21	1.42	1.00
1:B:335:PHE:HD2	3:B:783:HOH:O	1.58	0.85
1:B:594:VAL:HG23	1:B:595:THR:HG22	1.62	0.79
1:B:347:PRO:HG2	1:B:562:HIS:CE1	2.19	0.78
1:B:522:ALA:HB1	1:B:550:MET:HE3	1.72	0.71
1:A:515:LEU:HD22	1:A:566:ILE:HG13	1.74	0.69
1:B:562:HIS:HB3	1:B:567:TYR:CE2	2.27	0.68
1:A:466:ALA:HB1	1:A:514:VAL:HG23	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:355:LEU:HD23	1:B:396:TYR:OH	1.95	0.67
1:B:327:LEU:N	1:B:609:THR:O	2.31	0.64
1:B:336:ARG:HG3	3:B:706:HOH:O	1.98	0.64
1:B:335:PHE:CD2	3:B:783:HOH:O	2.41	0.62
1:B:351:THR:HG22	1:B:352:TRP:H	1.65	0.60
1:B:562:HIS:HB3	1:B:567:TYR:HE2	1.67	0.60
1:B:347:PRO:HG2	1:B:562:HIS:ND1	2.17	0.59
1:B:548:ALA:HB2	1:B:589:ASP:CG	2.28	0.59
1:B:454:ALA:HB3	1:B:490:TYR:HE1	1.67	0.58
1:B:472:LEU:HB3	1:B:490:TYR:HB3	1.87	0.56
1:B:566:ILE:HB	1:B:584:TYR:HB3	1.87	0.56
1:B:441:GLU:HB3	1:B:452:LEU:HD23	1.87	0.55
1:A:416:ILE:HD11	1:A:427:ALA:HB1	1.90	0.54
1:B:352:TRP:HE1	1:B:595:THR:HG1	1.48	0.54
1:B:397:ASN:OD1	1:B:399:MET:HB2	2.08	0.54
1:B:354:ARG:HD2	3:B:701:HOH:O	2.08	0.53
1:B:455:PRO:O	1:B:497:TRP:CD1	2.62	0.53
1:B:466:ALA:HB1	1:B:514:VAL:HG23	1.92	0.52
1:A:550:MET:HE3	1:A:568:VAL:CG2	2.29	0.52
1:B:592:SER:OG	1:B:594:VAL:HG13	2.10	0.51
1:B:447:ARG:O	1:B:449:GLU:HG3	2.10	0.51
1:B:519:ILE:O	1:B:536:ARG:HA	2.11	0.50
1:B:352:TRP:NE1	1:B:595:THR:OG1	2.26	0.50
1:B:584:TYR:HB2	1:B:591:TRP:CZ3	2.46	0.50
1:A:566:ILE:HB	1:A:584:TYR:HB3	1.95	0.48
1:A:426:TYR:CZ	1:A:442:ARG:HD3	2.49	0.48
1:B:421:ILE:HD11	1:B:472:LEU:HB2	1.95	0.48
1:A:346:ASN:CG	1:A:349:ASP:HB2	2.39	0.47
1:A:562:HIS:HB3	1:A:567:TYR:CE2	2.50	0.47
1:B:356:ALA:HB2	1:B:401:ASN:ND2	2.28	0.47
1:A:373:LEU:HB3	1:A:395:CYS:SG	2.55	0.47
1:B:550:MET:HE2	1:B:568:VAL:HG21	1.97	0.47
1:B:454:ALA:HB2	1:B:495:ASN:ND2	2.30	0.47
1:A:550:MET:CE	1:A:568:VAL:HG21	2.30	0.46
1:A:557:LEU:HD23	1:A:557:LEU:H	1.80	0.46
1:B:395:CYS:O	1:B:403:TRP:HA	2.16	0.46
1:B:347:PRO:CG	1:B:562:HIS:CE1	2.95	0.45
1:B:443:TYR:HB2	1:B:450:TRP:CE2	2.52	0.45
1:B:447:ARG:NH1	1:B:447:ARG:CG	2.75	0.45
1:B:550:MET:CE	1:B:568:VAL:HG11	2.47	0.45
1:B:454:ALA:HB3	1:B:490:TYR:CE1	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:490:TYR:OH	1:B:495:ASN:OD1	2.24	0.44
1:A:362:ARG:NH1	1:A:394:ASP:OD1	2.50	0.43
1:B:334:TYR:CG	2:C:82:GLU:HG2	2.53	0.43
1:A:472:LEU:HB3	1:A:490:TYR:HB3	1.99	0.43
1:B:335:PHE:CD1	1:B:600:GLY:O	2.71	0.43
1:A:466:ALA:HB2	1:A:512:VAL:HG12	2.00	0.43
1:B:418:VAL:HA	1:B:426:TYR:O	2.18	0.43
1:A:490:TYR:HB2	1:A:497:TRP:CE2	2.54	0.43
1:B:431:SER:HB3	1:B:461:ILE:HG21	2.01	0.43
1:B:548:ALA:HB2	1:B:589:ASP:OD1	2.19	0.43
1:B:554:ARG:HG3	1:B:557:LEU:HD22	2.01	0.43
1:B:352:TRP:CZ2	1:B:597:MET:HA	2.54	0.43
1:A:512:VAL:HA	1:A:520:TYR:O	2.19	0.42
1:B:351:THR:HG22	1:B:352:TRP:N	2.30	0.42
1:B:562:HIS:O	1:B:563:GLN:C	2.61	0.42
1:A:365:LEU:HD12	1:A:376:ALA:HB1	2.02	0.42
1:B:530:GLN:HG2	1:B:573:ASP:HA	2.01	0.42
1:A:338:SER:HB3	1:A:361:PRO:C	2.45	0.42
1:B:595:THR:OG1	1:B:596:ARG:N	2.53	0.42
1:B:365:LEU:H	1:B:365:LEU:HD23	1.85	0.41
1:B:349:ASP:OD2	1:B:349:ASP:N	2.53	0.41
1:B:336:ARG:N	3:B:706:HOH:O	2.54	0.41
1:B:579:ASP:HB2	1:B:596:ARG:C	2.45	0.41
1:B:416:ILE:HD11	1:B:427:ALA:HB1	2.02	0.41
1:B:443:TYR:HB2	1:B:450:TRP:CD2	2.56	0.41
1:B:553:ARG:H	1:B:553:ARG:HG3	1.63	0.41
1:B:345:TYR:HB2	1:B:352:TRP:CZ3	2.56	0.41
1:B:425:ILE:HD12	1:B:443:TYR:CD2	2.55	0.40
1:A:338:SER:O	1:A:361:PRO:HB3	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:517:ASN:OD1	1:B:470:ARG:NH1[2_555]	1.55	0.65

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	281/336 (84%)	271 (96%)	10 (4%)	0	100	100
1	B	281/336 (84%)	262 (93%)	17 (6%)	2 (1%)	18	14
2	C	7/9 (78%)	7 (100%)	0	0	100	100
All	All	569/681 (84%)	540 (95%)	27 (5%)	2 (0%)	30	28

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	597	MET
1	B	598	THR

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	228/275 (83%)	227 (100%)	1 (0%)	84	89
1	B	228/275 (83%)	222 (97%)	6 (3%)	40	45
2	C	5/5 (100%)	5 (100%)	0	100	100
All	All	461/555 (83%)	454 (98%)	7 (2%)	57	64

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	517	ASN

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Mol	Chain	Res	Type
1	B	336	ARG
1	B	451	HIS
1	B	590	THR
1	B	591	TRP
1	B	592	SER
1	B	596	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	517	ASN
1	A	552	HIS
1	B	401	ASN
1	B	414	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 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	283/336 (84%)	0.26	21 (7%) 20 22	20, 28, 50, 76	0
1	B	283/336 (84%)	1.67	106 (37%) 1 1	19, 42, 80, 95	0
2	C	7/9 (77%)	1.11	2 (28%) 1 1	29, 32, 51, 53	0
All	All	573/681 (84%)	0.97	129 (22%) 2 2	19, 33, 70, 95	0

All (129) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	335	PHE	8.0
1	B	594	VAL	7.0
1	B	590	THR	6.9
1	B	329	TYR	6.2
1	B	327	LEU	5.7
1	A	338	SER	5.6
1	B	592	SER	5.6
1	B	350	GLY	5.5
1	B	591	TRP	5.4
1	B	345	TYR	5.1
1	A	337	GLN	5.0
1	B	454	ALA	4.9
1	A	335	PHE	4.8
1	B	344	ALA	4.7
1	B	424	HIS	4.7
1	B	328	ILE	4.7
1	B	395	CYS	4.6
1	B	352	TRP	4.5
1	B	609	THR	4.4
1	B	548	ALA	4.4
1	B	396	TYR	4.3
1	B	351	THR	4.3
1	B	608	VAL	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	567	TYR	4.2
1	B	588	THR	4.2
1	B	423	GLY	4.2
1	B	370	VAL	4.2
1	B	336	ARG	4.1
1	B	398	PRO	4.1
1	B	353	LEU	4.0
1	B	562	HIS	4.0
1	B	375	TYR	3.9
1	B	421	ILE	3.9
1	B	589	ASP	3.9
1	B	451	HIS	3.8
1	B	445	PRO	3.8
1	B	561	VAL	3.8
1	B	399	MET	3.8
1	B	584	TYR	3.7
1	B	425	ILE	3.7
1	B	397	ASN	3.6
1	B	581	VAL	3.6
1	B	405	PRO	3.6
1	B	583	CYS	3.6
1	A	336	ARG	3.5
1	B	587	ASP	3.5
1	B	470	ARG	3.5
1	B	595	THR	3.5
1	B	493	GLU	3.5
1	B	369	VAL	3.5
1	B	593	GLU	3.4
1	B	443	TYR	3.4
1	A	547	VAL	3.4
1	B	491	TYR	3.4
2	C	76	GLY	3.4
1	B	373	LEU	3.4
1	B	597	MET	3.3
1	B	347	PRO	3.3
1	B	564	GLY	3.3
1	B	514	VAL	3.3
1	B	566	ILE	3.3
1	B	346	ASN	3.3
1	B	420	VAL	3.2
1	B	342	LEU	3.2
1	B	518	CYS	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	407	ALA	3.2
1	A	384	PRO	3.2
1	B	376	ALA	3.1
1	B	607	ALA	3.1
1	B	446	GLU	3.1
1	B	578	LEU	3.1
1	A	387	ASN	3.1
1	B	422	ASP	3.1
1	B	596	ARG	3.0
1	B	453	VAL	3.0
1	B	330	THR	3.0
1	B	452	LEU	2.9
1	B	471	LEU	2.9
1	B	442	ARG	2.9
1	B	374	LEU	2.9
1	B	371	GLY	2.9
1	B	494	ARG	2.9
1	B	496	GLU	2.9
1	B	348	SER	2.9
1	A	349	ASP	2.8
1	B	360	VAL	2.8
2	C	78	PRO	2.8
1	B	349	ASP	2.8
1	A	386	GLY	2.8
1	B	516	HIS	2.7
1	B	449	GLU	2.7
1	A	470	ARG	2.7
1	A	446	GLU	2.7
1	B	563	GLN	2.6
1	B	586	PRO	2.6
1	A	385	ASP	2.6
1	B	450	TRP	2.6
1	B	492	PRO	2.6
1	B	400	THR	2.6
1	B	403	TRP	2.6
1	A	494	ARG	2.6
1	B	468	LEU	2.6
1	A	546	PHE	2.5
1	B	585	ASP	2.5
1	B	549	PRO	2.5
1	B	598	THR	2.5
1	A	452	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	393	LEU	2.5
1	A	491	TYR	2.4
1	B	447	ARG	2.4
1	A	454	ALA	2.4
1	B	600	GLY	2.3
1	B	513	CYS	2.3
1	B	402	GLN	2.3
1	A	498	ARG	2.3
1	B	355	LEU	2.3
1	A	399	MET	2.2
1	B	553	ARG	2.2
1	B	467	VAL	2.2
1	B	368	CYS	2.2
1	A	493	GLU	2.2
1	B	495	ASN	2.2
1	B	582	GLU	2.1
1	B	575	HIS	2.1
1	A	492	PRO	2.1
1	B	552	HIS	2.1
1	B	565	ARG	2.1
1	B	568	VAL	2.0
1	B	359	GLN	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.