



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:55 AM UTC

PDB ID : 6KGJ / pdb_00006kgj
Title : M1Q-hNTAQ1 C28S
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Deposited on : 2019-07-11
Resolution : 1.80 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

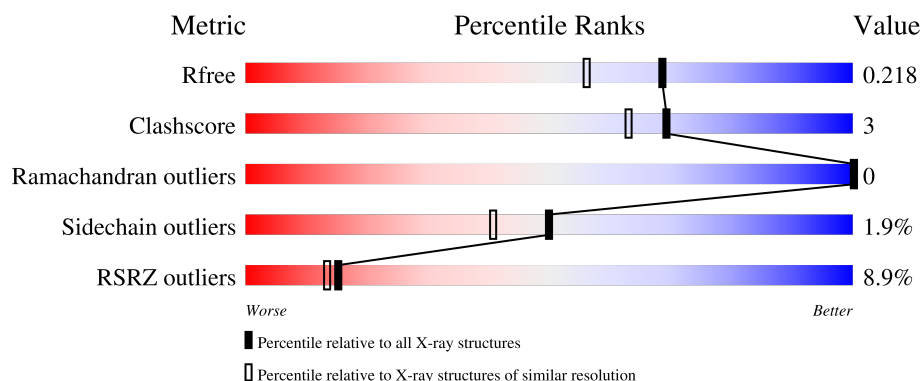
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	202	10% 42% 52% 5%
1	B	202	7% 40% 54% 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3426 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 Protein N-terminal glutamine amidohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	202	Total	C	N	O	S	0	0	0
			1635	1042	273	309	11			
1	A	202	Total	C	N	O	S	0	0	0
			1635	1042	273	309	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	GLN	-	expression tag	UNP Q96HA8
B	28	SER	CYS	engineered mutation	UNP Q96HA8
B	32	VAL	ILE	variant	UNP Q96HA8
B	93	SER	ASN	variant	UNP Q96HA8
B	116	ILE	PHE	variant	UNP Q96HA8
B	134	CYS	ARG	variant	UNP Q96HA8
A	1	GLN	-	expression tag	UNP Q96HA8
A	28	SER	CYS	engineered mutation	UNP Q96HA8
A	32	VAL	ILE	variant	UNP Q96HA8
A	93	SER	ASN	variant	UNP Q96HA8
A	116	ILE	PHE	variant	UNP Q96HA8
A	134	CYS	ARG	variant	UNP Q96HA8

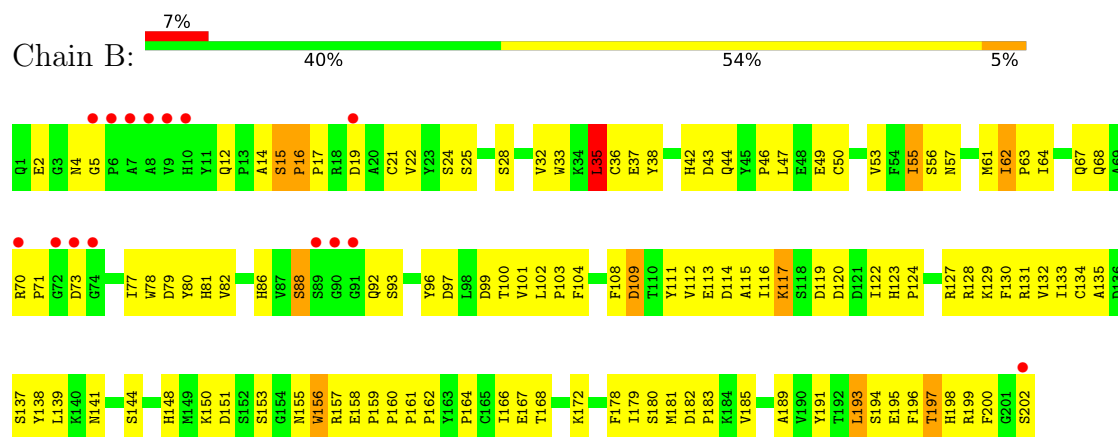
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	75	Total	O	0	0
			75	75		
2	A	81	Total	O	0	0
			81	81		

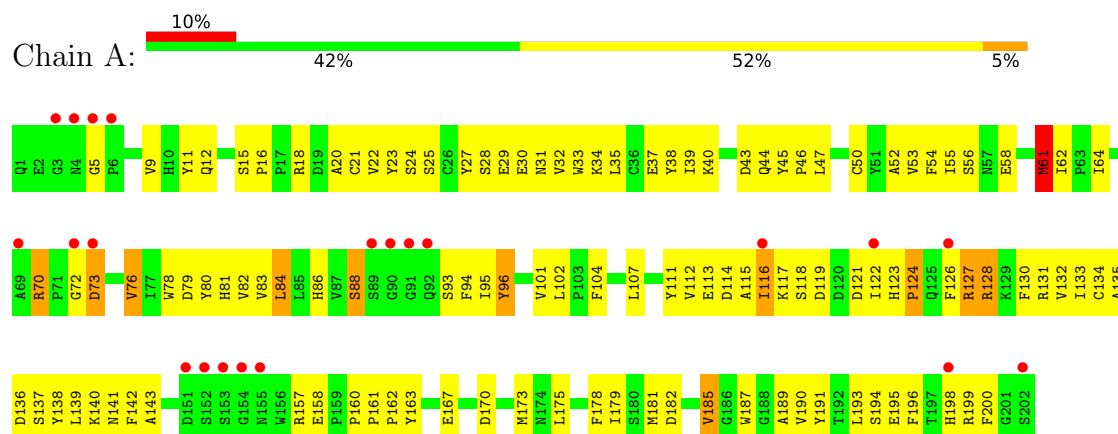
3 Residue-property plots [i](#)

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: Protein N-terminal glutamine amidohydrolase



- Molecule 1: Protein N-terminal glutamine amidohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	95.64Å 105.78Å 140.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.47 – 1.80 35.47 – 1.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (35.47-1.80) 99.5 (35.47-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 1.81Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.211 , 0.223 0.217 , 0.218	Depositor DCC
R_{free} test set	2000 reflections (3.05%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 27.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3426	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.54% 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	2.30	110/1690 (6.5%)	1.19	13/2300 (0.6%)
1	B	2.28	104/1690 (6.2%)	1.20	12/2300 (0.5%)
All	All	2.29	214/3380 (6.3%)	1.20	25/4600 (0.5%)

All (214) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	161	PRO	C-O	-11.06	1.15	1.25
1	A	15	SER	C-O	-10.05	1.15	1.24
1	B	62	ILE	C-O	-9.48	1.15	1.23
1	A	133	ILE	C-O	-9.43	1.14	1.24
1	A	185	VAL	C-O	-8.38	1.15	1.24
1	A	107	LEU	C-O	-8.24	1.14	1.23
1	B	82	VAL	C-O	-8.19	1.15	1.24
1	B	104	PHE	C-O	-8.17	1.17	1.25
1	A	111	TYR	C-O	-8.13	1.14	1.24
1	B	15	SER	C-O	-8.11	1.17	1.24
1	A	158	GLU	C-O	-8.01	1.15	1.24
1	A	130	PHE	C-O	-7.93	1.15	1.24
1	A	127	ARG	C-O	-7.79	1.14	1.23
1	B	78	TRP	C-O	-7.73	1.14	1.23
1	A	81	HIS	C-O	-7.73	1.14	1.23
1	B	64	ILE	C-O	-7.68	1.16	1.24
1	A	160	PRO	C-O	-7.66	1.16	1.24
1	A	117	LYS	C-O	-7.57	1.16	1.24
1	A	193	LEU	C-O	-7.55	1.15	1.24
1	B	191	TYR	C-O	-7.50	1.14	1.23
1	A	45	TYR	C-O	-7.50	1.16	1.24
1	A	132	VAL	C-O	-7.43	1.16	1.24
1	A	53	VAL	C-O	-7.43	1.15	1.24
1	A	35	LEU	C-O	-7.39	1.15	1.24
1	B	2	GLU	C-O	-7.38	1.15	1.23
1	A	82	VAL	C-O	-7.37	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	133	ILE	C-O	-7.35	1.16	1.24
1	A	95	ILE	C-O	-7.32	1.16	1.24
1	B	22	VAL	C-O	-7.28	1.16	1.24
1	A	32	VAL	C-O	-7.22	1.15	1.24
1	B	53	VAL	C-O	-7.19	1.16	1.24
1	B	57	ASN	C-O	-7.17	1.16	1.23
1	A	86	HIS	C-O	-7.15	1.15	1.24
1	A	112	VAL	C-O	-7.12	1.15	1.24
1	B	117	LYS	C-O	-7.12	1.15	1.24
1	A	157	ARG	C-O	-7.10	1.15	1.24
1	A	29	GLU	C-O	-7.09	1.15	1.24
1	A	94	PHE	C-O	-7.07	1.15	1.23
1	B	63	PRO	C-O	-7.00	1.16	1.23
1	A	9	VAL	C-O	-6.96	1.16	1.24
1	A	25	SER	C-O	-6.95	1.15	1.23
1	A	58	GLU	C-O	-6.95	1.16	1.24
1	A	138	TYR	C-O	-6.94	1.16	1.24
1	A	21	CYS	C-O	-6.93	1.15	1.23
1	B	80	TYR	C-O	-6.92	1.15	1.23
1	A	24	SER	C-O	-6.91	1.16	1.24
1	B	139	LEU	C-O	-6.85	1.16	1.24
1	A	195	GLU	C-O	-6.84	1.16	1.24
1	A	28	SER	C-O	-6.81	1.15	1.24
1	B	138	TYR	C-O	-6.77	1.16	1.24
1	B	178	PHE	C-O	-6.76	1.15	1.24
1	B	180	SER	C-O	-6.76	1.15	1.23
1	A	54	PHE	C-O	-6.73	1.15	1.23
1	A	46	PRO	C-O	-6.71	1.16	1.23
1	B	124	PRO	C-O	-6.70	1.15	1.24
1	A	135	ALA	C-O	-6.68	1.16	1.24
1	A	140	LYS	C-O	-6.67	1.16	1.24
1	B	195	GLU	C-O	-6.64	1.16	1.24
1	A	142	PHE	C-O	-6.63	1.16	1.23
1	B	182	ASP	C-O	-6.62	1.15	1.24
1	A	104	PHE	C-O	-6.61	1.19	1.25
1	B	158	GLU	C-O	-6.60	1.17	1.24
1	B	131	ARG	C-O	-6.60	1.16	1.24
1	B	159	PRO	C-O	-6.58	1.17	1.24
1	B	197	THR	C-O	-6.53	1.16	1.24
1	B	189	ALA	C-O	-6.53	1.15	1.23
1	B	122	ILE	C-O	-6.52	1.15	1.24
1	A	118	SER	C-O	-6.50	1.16	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	191	TYR	C-O	-6.49	1.16	1.23
1	A	12	GLN	C-O	-6.45	1.15	1.24
1	A	143	ALA	C-O	-6.45	1.16	1.23
1	A	134	CYS	C-O	-6.44	1.16	1.23
1	B	111	TYR	C-O	-6.44	1.16	1.24
1	B	123	HIS	C-O	-6.43	1.15	1.23
1	B	127	ARG	C-O	-6.41	1.16	1.23
1	B	135	ALA	C-O	-6.41	1.16	1.24
1	A	37	GLU	C-O	-6.41	1.16	1.24
1	A	38	TYR	C-O	-6.41	1.16	1.24
1	A	62	ILE	C-O	-6.41	1.18	1.23
1	A	162	PRO	C-O	-6.40	1.17	1.23
1	A	189	ALA	C-O	-6.36	1.15	1.23
1	A	170	ASP	C-O	-6.34	1.16	1.24
1	B	93	SER	C-O	-6.33	1.16	1.23
1	B	81	HIS	C-O	-6.32	1.16	1.23
1	A	161	PRO	C-O	-6.24	1.17	1.24
1	B	130	PHE	C-O	-6.23	1.16	1.24
1	A	137	SER	C-O	-6.22	1.16	1.24
1	A	198	HIS	C-O	-6.21	1.16	1.24
1	B	113	GLU	C-O	-6.20	1.17	1.24
1	B	71	PRO	C-O	-6.20	1.16	1.23
1	B	96	TYR	C-O	-6.18	1.16	1.24
1	A	141	ASN	C-O	-6.18	1.15	1.23
1	B	167	GLU	C-O	-6.16	1.16	1.23
1	A	22	VAL	C-O	-6.14	1.17	1.24
1	B	55	ILE	C-O	-6.14	1.17	1.24
1	A	44	GLN	C-O	-6.14	1.17	1.24
1	A	178	PHE	C-O	-6.13	1.15	1.23
1	B	79	ASP	C-O	-6.12	1.16	1.24
1	A	93	SER	C-O	-6.11	1.16	1.23
1	B	198	HIS	C-O	-6.08	1.17	1.24
1	A	27	TYR	C-O	-6.08	1.16	1.23
1	B	200	PHE	C-O	-6.08	1.15	1.24
1	A	16	PRO	C-O	-6.06	1.17	1.24
1	A	88	SER	C-O	-6.05	1.15	1.23
1	A	56	SER	C-O	-6.05	1.18	1.24
1	A	114	ASP	C-O	-6.05	1.17	1.24
1	A	84	LEU	C-O	-6.03	1.16	1.24
1	A	179	ILE	C-O	-6.02	1.17	1.24
1	A	102	LEU	C-O	-6.01	1.16	1.23
1	B	164	PRO	C-O	-6.00	1.16	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	156	TRP	C-O	-6.00	1.16	1.23
1	A	128	ARG	C-O	-5.99	1.16	1.24
1	B	128	ARG	C-O	-5.98	1.16	1.24
1	A	199	ARG	C-O	-5.96	1.17	1.24
1	A	34	LYS	C-O	-5.94	1.17	1.24
1	B	56	SER	C-O	-5.92	1.18	1.24
1	B	166	ILE	C-O	-5.91	1.17	1.24
1	B	46	PRO	C-O	-5.90	1.16	1.23
1	B	153	SER	C-O	-5.90	1.16	1.24
1	B	35	LEU	C-O	-5.88	1.17	1.24
1	B	137	SER	C-O	-5.88	1.17	1.24
1	A	61	MET	C-O	-5.87	1.16	1.23
1	A	40	LYS	C-O	-5.87	1.17	1.24
1	A	18	ARG	C-O	-5.85	1.17	1.24
1	A	163	TYR	C-O	-5.84	1.18	1.24
1	A	83	VAL	C-O	-5.83	1.16	1.23
1	A	30	GLU	C-O	-5.81	1.17	1.24
1	B	86	HIS	C-O	-5.81	1.17	1.24
1	B	120	ASP	C-O	-5.80	1.17	1.24
1	B	157	ARG	C-O	-5.80	1.17	1.24
1	B	144	SER	C-O	-5.79	1.17	1.23
1	A	31	ASN	C-O	-5.79	1.17	1.24
1	B	108	PHE	C-O	-5.77	1.17	1.24
1	B	61	MET	C-O	-5.76	1.16	1.23
1	A	139	LEU	C-O	-5.76	1.17	1.24
1	B	134	CYS	C-O	-5.74	1.17	1.23
1	A	39	ILE	C-O	-5.74	1.17	1.24
1	B	109	ASP	C-O	-5.73	1.17	1.24
1	A	79	ASP	C-O	-5.73	1.17	1.24
1	B	185	VAL	C-O	-5.70	1.17	1.24
1	A	182	ASP	C-O	-5.70	1.17	1.24
1	B	100	THR	C-O	-5.70	1.16	1.23
1	A	161	PRO	C-N	-5.68	1.29	1.33
1	B	196	PHE	C-O	-5.67	1.17	1.24
1	B	101	VAL	C-O	-5.66	1.18	1.24
1	B	194	SER	C-O	-5.66	1.17	1.24
1	B	162	PRO	CA-CB	-5.66	1.50	1.53
1	B	32	VAL	C-O	-5.60	1.17	1.24
1	B	193	LEU	C-O	-5.59	1.17	1.24
1	B	119	ASP	C-O	-5.58	1.16	1.24
1	A	20	ALA	C-O	-5.58	1.16	1.23
1	B	71	PRO	CA-C	-5.55	1.47	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	80	TYR	C-O	-5.55	1.17	1.23
1	A	173	MET	C-O	-5.55	1.17	1.23
1	A	55	ILE	C-O	-5.55	1.18	1.24
1	B	67	GLN	C-O	-5.55	1.17	1.24
1	A	101	VAL	C-O	-5.55	1.18	1.24
1	B	12	GLN	C-O	-5.55	1.17	1.24
1	B	33	TRP	C-O	-5.54	1.17	1.24
1	B	168	THR	C-O	-5.52	1.16	1.23
1	B	199	ARG	C-O	-5.52	1.17	1.24
1	A	122	ILE	C-O	-5.51	1.17	1.23
1	B	115	ALA	C-O	-5.50	1.17	1.24
1	B	160	PRO	C-O	-5.48	1.18	1.24
1	A	175	LEU	C-O	-5.47	1.17	1.24
1	B	183	PRO	C-O	-5.46	1.17	1.24
1	A	33	TRP	C-O	-5.45	1.17	1.24
1	B	25	SER	C-O	-5.45	1.17	1.23
1	A	50	CYS	C-O	-5.44	1.17	1.23
1	A	124	PRO	C-O	-5.42	1.17	1.24
1	B	97	ASP	C-O	-5.41	1.17	1.24
1	B	50	CYS	C-O	-5.40	1.17	1.23
1	B	21	CYS	C-O	-5.38	1.17	1.23
1	A	47	LEU	C-O	-5.37	1.17	1.24
1	A	181	MET	C-O	-5.37	1.16	1.24
1	B	132	VAL	C-O	-5.35	1.18	1.24
1	B	114	ASP	C-O	-5.34	1.17	1.24
1	B	179	ILE	C-O	-5.31	1.17	1.24
1	B	112	VAL	C-O	-5.29	1.17	1.24
1	A	196	PHE	C-O	-5.29	1.18	1.24
1	B	155	ASN	C-O	-5.28	1.17	1.23
1	A	190	VAL	C-O	-5.27	1.18	1.24
1	A	113	GLU	C-O	-5.27	1.18	1.24
1	B	77	ILE	C-O	-5.27	1.18	1.24
1	A	131	ARG	C-O	-5.25	1.17	1.24
1	A	78	TRP	C-O	-5.25	1.17	1.23
1	A	64	ILE	C-O	-5.24	1.18	1.24
1	B	17	PRO	C-O	-5.23	1.17	1.23
1	A	11	TYR	C-O	-5.22	1.17	1.23
1	B	129	LYS	C-O	-5.21	1.17	1.23
1	B	148	HIS	C-O	-5.20	1.17	1.24
1	B	24	SER	C-O	-5.20	1.18	1.24
1	B	151	ASP	C-O	-5.18	1.18	1.23
1	A	200	PHE	C-O	-5.16	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	47	LEU	C-O	-5.15	1.17	1.24
1	B	181	MET	C-O	-5.12	1.17	1.24
1	A	116	ILE	C-O	-5.12	1.18	1.24
1	B	49	GLU	C-O	-5.11	1.17	1.24
1	A	43	ASP	C-O	-5.11	1.17	1.24
1	B	14	ALA	C-O	-5.10	1.17	1.24
1	A	121	ASP	C-O	-5.10	1.17	1.24
1	A	167	GLU	C-O	-5.09	1.17	1.23
1	B	36	CYS	C-O	-5.07	1.18	1.24
1	B	19	ASP	N-CA	5.07	1.52	1.46
1	A	52	ALA	C-O	-5.06	1.17	1.23
1	A	76	VAL	N-CA	-5.06	1.40	1.46
1	A	187	TRP	C-O	-5.05	1.17	1.23
1	B	37	GLU	C-O	-5.05	1.18	1.24
1	B	43	ASP	CA-C	-5.05	1.46	1.52
1	B	28	SER	C-O	-5.04	1.17	1.24
1	A	96	TYR	C-O	-5.04	1.18	1.24
1	A	136	ASP	C-O	-5.04	1.18	1.24
1	B	172	LYS	C-O	-5.04	1.17	1.24
1	A	194	SER	C-O	-5.02	1.18	1.24

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	73	ASP	N-CA-C	12.81	138.09	110.80
1	B	141	ASN	N-CA-C	6.93	122.41	113.88
1	B	114	ASP	N-CA-C	6.43	119.25	111.40
1	B	115	ALA	N-CA-C	6.26	118.19	111.36
1	B	4	ASN	N-CA-C	6.16	123.93	110.80
1	A	141	ASN	N-CA-C	6.05	121.32	113.88
1	A	115	ALA	N-CA-C	6.04	117.94	111.36
1	B	5	GLY	N-CA-C	-5.96	100.19	112.34
1	A	72	GLY	N-CA-C	-5.75	99.56	113.18
1	B	199	ARG	N-CA-C	5.73	117.33	111.14
1	B	19	ASP	N-CA-C	5.73	119.37	112.38
1	B	44	GLN	N-CA-C	5.70	117.30	111.14
1	A	5	GLY	CA-C-N	5.63	125.58	119.78
1	A	5	GLY	C-N-CA	5.63	125.58	119.78
1	B	4	ASN	CA-C-N	-5.50	113.23	121.87
1	B	4	ASN	C-N-CA	-5.50	113.23	121.87
1	A	114	ASP	N-CA-C	5.46	117.31	111.36
1	A	70	ARG	CA-C-N	5.36	125.28	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	ARG	C-N-CA	5.36	125.28	119.76
1	A	23	TYR	N-CA-C	5.31	117.15	108.55
1	B	88	SER	N-CA-C	5.30	118.33	110.48
1	A	123	HIS	CA-C-N	5.22	125.03	119.28
1	A	123	HIS	C-N-CA	5.22	125.03	119.28
1	A	93	SER	CB-CA-C	5.07	119.31	109.33
1	B	16	PRO	N-CA-C	-5.02	104.58	110.70

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	1635	0	1530	7	0
1	B	1635	0	1530	15	0
2	A	81	0	0	0	0
2	B	75	0	0	0	0
All	All	3426	0	3060	22	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 3.

All (22) 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:70:ARG:HG3	1:A:76:VAL:HG22	1.72	0.71
1:A:61:MET:HE2	1:A:126:PHE:HE1	1.59	0.66
1:B:38:TYR:CE2	1:B:42:HIS:CE1	2.87	0.63
1:B:116:ILE:HG21	1:B:193:LEU:HD11	1.80	0.63
1:A:61:MET:HE2	1:A:126:PHE:CE1	2.36	0.61
1:B:193:LEU:O	1:B:197:THR:HG23	2.01	0.60
1:A:119:ASP:CG	1:A:127:ARG:HG2	2.29	0.58
1:B:116:ILE:HG21	1:B:193:LEU:CD1	2.35	0.56
1:A:116:ILE:HG23	1:A:128:ARG:CZ	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:LYS:HE2	1:B:156:TRP:CE2	2.48	0.48
1:B:150:LYS:HE2	1:B:156:TRP:CZ2	2.49	0.47
1:A:124:PRO:HA	1:A:127:ARG:HD2	1.95	0.47
1:B:15:SER:HB3	1:B:35:LEU:HD21	1.96	0.46
1:B:70:ARG:HE	1:B:73:ASP:CB	2.27	0.45
1:B:102:LEU:HB3	1:B:103:PRO:HD2	1.98	0.45
1:B:88:SER:HB3	1:B:92:GLN:O	2.18	0.43
1:B:15:SER:HB2	1:B:16:PRO:HD2	2.01	0.42
1:A:84:LEU:HB3	1:A:96:TYR:HB2	2.00	0.42
1:B:109:ASP:OD2	1:B:202:SER:HB2	2.19	0.42
1:B:102:LEU:HB3	1:B:103:PRO:CD	2.50	0.42
1:B:68:GLN:HG3	1:B:99:ASP:O	2.20	0.41
1:B:116:ILE:HG22	1:B:117:LYS:O	2.20	0.41

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	A	200/202 (99%)	196 (98%)	4 (2%)	0	100	100
1	B	200/202 (99%)	193 (96%)	7 (4%)	0	100	100
All	All	400/404 (99%)	389 (97%)	11 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	181/181 (100%)	177 (98%)	4 (2%)	45	34
1	B	181/181 (100%)	178 (98%)	3 (2%)	53	45
All	All	362/362 (100%)	355 (98%)	7 (2%)	50	41

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

Mol	Chain	Res	Type
1	B	35	LEU
1	B	55	ILE
1	B	62	ILE
1	A	61	MET
1	A	73	ASP
1	A	88	SER
1	A	185	VAL

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

Mol	Chain	Res	Type
1	B	42	HIS
1	B	92	GLN
1	B	155	ASN
1	A	1	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

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	A	202/202 (100%)	0.36	21 (10%)	11 10	16, 24, 46, 59	0
1	B	202/202 (100%)	0.39	15 (7%)	20 19	16, 25, 44, 55	0
All	All	404/404 (100%)	0.38	36 (8%)	15 13	16, 24, 45, 59	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	6	PRO	4.6
1	B	9	VAL	4.5
1	B	5	GLY	4.2
1	B	89	SER	3.9
1	B	72	GLY	3.9
1	B	10	HIS	3.6
1	A	5	GLY	3.5
1	A	152	SER	3.5
1	B	73	ASP	3.5
1	A	69	ALA	3.3
1	A	90	GLY	3.3
1	B	74	GLY	3.2
1	B	7	ALA	3.2
1	B	202	SER	3.1
1	A	116	ILE	3.1
1	A	122	ILE	2.9
1	B	90	GLY	2.9
1	A	89	SER	2.9
1	A	3	GLY	2.9
1	A	4	ASN	2.8
1	A	72	GLY	2.7
1	B	19	ASP	2.6
1	A	126	PHE	2.6
1	A	73	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	8	ALA	2.5
1	A	155	ASN	2.5
1	B	70	ARG	2.5
1	B	6	PRO	2.4
1	A	202	SER	2.4
1	A	198	HIS	2.3
1	A	91	GLY	2.2
1	A	92	GLN	2.2
1	A	153	SER	2.2
1	B	91	GLY	2.2
1	A	154	GLY	2.2
1	A	151	ASP	2.1

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.