



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 11:02 AM UTC

PDB ID : 7T0H / pdb_00007t0h
Title : Crystal structure of S25-39 Fab Unliganded 2
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Deposited on : 2021-11-29
Resolution : 1.30 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

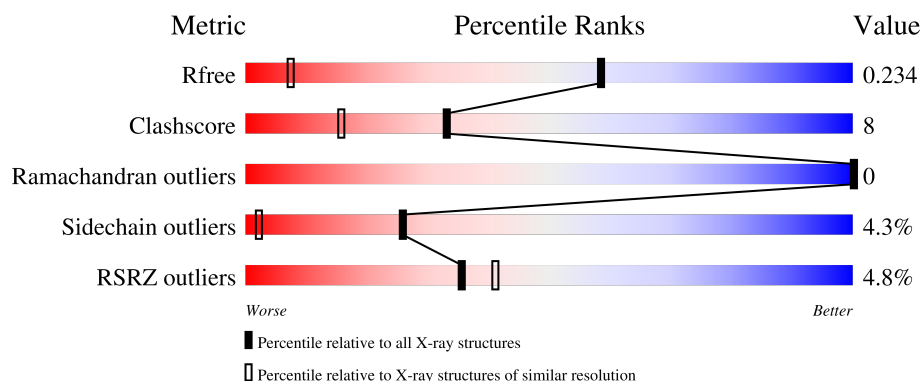
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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	1553 (1.30-1.30)
Clashscore	190562	1595 (1.30-1.30)
Ramachandran outliers	187476	1551 (1.30-1.30)
Sidechain outliers	187428	1551 (1.30-1.30)
RSRZ outliers	180081	1549 (1.30-1.30)

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	H	222	
2	L	219	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BR	H	301	-	-	X	-
3	BR	H	302	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4089 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 S25-39 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	215	Total	C	N	O	S	0	12	0
			1731	1093	287	339	12			

- Molecule 2 is a protein called S25-39 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	219	Total	C	N	O	S	0	10	0
			1761	1090	301	359	11			

- Molecule 3 is BROMIDE ION (CCD ID: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	3	Total	Br	0	0
			3	3		
3	L	2	Total	Br	0	0
			2	2		

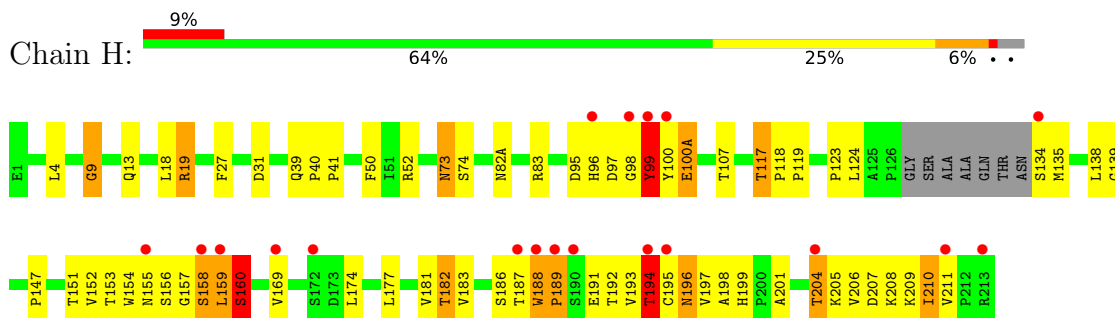
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	318	Total	O	0	1
			319	319		
4	L	273	Total	O	0	0
			273	273		

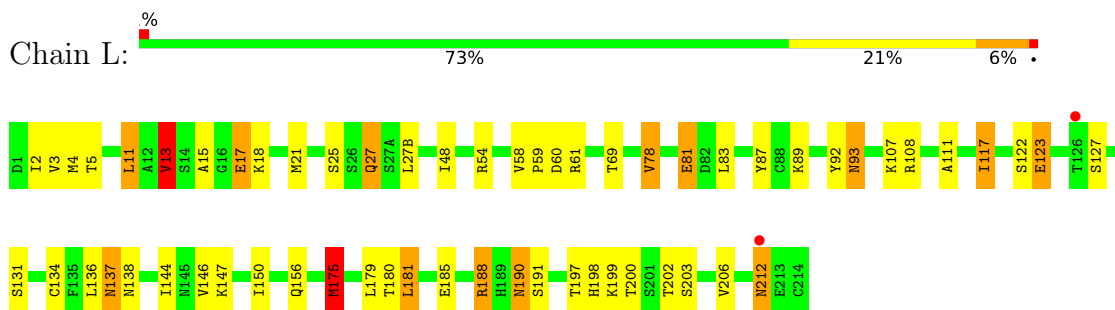
3 Residue-property plots [i](#)

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: S25-39 Fab heavy chain



- Molecule 2: S25-39 Fab light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.59Å 63.13Å 61.88Å 90.00° 99.36° 90.00°	Depositor
Resolution (Å)	40.00 – 1.30 40.00 – 1.30	Depositor EDS
% Data completeness (in resolution range)	91.4 (40.00-1.30) 91.7 (40.00-1.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.85 (at 1.30Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.148 , 0.200 (Not available) , 0.234	Depositor DCC
R_{free} test set	4735 reflections (4.54%)	wwPDB-VP
Wilson B-factor (Å ²)	13.6	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4089	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.45% 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: BR

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	H	2.15	64/1776 (3.6%)	1.73	46/2423 (1.9%)
2	L	2.08	46/1799 (2.6%)	1.66	31/2435 (1.3%)
All	All	2.12	110/3575 (3.1%)	1.70	77/4858 (1.6%)

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	H	0	3
2	L	0	1
All	All	0	4

All (110) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	21	MET	SD-CE	-13.28	1.46	1.79
2	L	137	ASN	CG-ND2	-11.71	1.08	1.33
2	L	27	GLN	CD-OE1	11.18	1.44	1.23
2	L	4	MET	CG-SD	-10.69	1.54	1.80
1	H	208	LYS	C-O	10.57	1.35	1.24
1	H	9	GLY	C-N	-9.25	1.23	1.33
1	H	98	GLY	N-CA	9.11	1.59	1.44
2	L	69	THR	CA-C	9.01	1.64	1.52
1	H	139	GLY	N-CA	8.81	1.54	1.45
1	H	96	HIS	C-O	-8.74	1.13	1.23
1	H	207	ASP	CA-C	8.49	1.62	1.52
2	L	5	THR	N-CA	8.39	1.56	1.45
1	H	99	TYR	N-CA	8.25	1.54	1.46
1	H	196	ASN	CA-CB	7.93	1.64	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	160	SER	N-CA	7.92	1.56	1.46
1	H	31	ASP	CG-OD2	7.84	1.40	1.25
1	H	174	LEU	C-O	-7.83	1.14	1.23
1	H	118	PRO	CA-C	7.67	1.59	1.52
2	L	78	VAL	CA-CB	-7.59	1.46	1.54
1	H	18	LEU	CB-CG	-7.50	1.38	1.53
1	H	206	VAL	C-O	7.30	1.31	1.24
1	H	192	THR	C-O	7.29	1.32	1.23
2	L	198	HIS	CA-C	7.28	1.61	1.52
1	H	153	THR	N-CA	7.27	1.56	1.46
1	H	193	VAL	CA-C	-7.16	1.43	1.52
2	L	27(B)	LEU	N-CA	7.11	1.54	1.46
1	H	209	LYS	C-O	7.10	1.32	1.23
1	H	160	SER	C-O	7.08	1.32	1.24
1	H	39	GLN	CA-CB	-7.07	1.46	1.53
2	L	11	LEU	CA-C	-7.00	1.43	1.52
2	L	11	LEU	CA-CB	-6.92	1.41	1.53
2	L	156	GLN	N-CA	6.92	1.54	1.46
1	H	123	PRO	CA-C	6.89	1.61	1.52
1	H	188	TRP	C-O	6.86	1.32	1.24
1	H	196	ASN	CA-C	6.82	1.61	1.52
2	L	15	ALA	C-O	6.75	1.31	1.23
1	H	158[A]	SER	CA-C	6.74	1.62	1.52
1	H	158[B]	SER	CA-C	6.74	1.62	1.52
2	L	188	ARG	CD-NE	6.74	1.55	1.46
2	L	191	SER	CA-CB	-6.73	1.43	1.53
1	H	82(A)	ASN	N-CA	6.71	1.53	1.45
1	H	197	VAL	C-O	-6.69	1.16	1.24
1	H	181	VAL	N-CA	-6.55	1.38	1.46
1	H	189	PRO	C-N	6.51	1.42	1.33
1	H	27	PHE	CA-C	-6.47	1.44	1.52
1	H	13	GLN	CD-OE1	6.45	1.35	1.23
2	L	111	ALA	CA-C	-6.39	1.44	1.52
1	H	204	THR	C-O	6.37	1.31	1.24
1	H	117	THR	CA-CB	6.27	1.64	1.53
1	H	196	ASN	CG-ND2	-6.23	1.20	1.33
1	H	39	GLN	C-N	-6.16	1.28	1.33
1	H	119	PRO	N-CA	6.16	1.54	1.47
1	H	152	VAL	C-O	-6.15	1.17	1.24
1	H	199	HIS	C-O	-6.15	1.17	1.24
2	L	181	LEU	C-O	-6.14	1.16	1.23
1	H	192	THR	CA-C	6.14	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	180	THR	C-O	6.14	1.31	1.24
2	L	60	ASP	CB-CG	-6.13	1.36	1.52
1	H	100(A)	GLU	CA-CB	6.12	1.61	1.52
2	L	27	GLN	CG-CD	6.08	1.67	1.52
2	L	202	THR	N-CA	-6.07	1.38	1.46
1	H	97	ASP	CG-OD1	6.06	1.36	1.25
1	H	211	VAL	C-N	6.05	1.41	1.33
1	H	194	THR	CA-C	6.05	1.59	1.52
1	H	153	THR	CA-C	6.01	1.61	1.52
1	H	186	SER	C-O	-6.01	1.16	1.24
1	H	210	ILE	N-CA	5.99	1.53	1.46
1	H	19	ARG	CD-NE	-5.93	1.38	1.46
2	L	61	ARG	CZ-NH2	5.92	1.41	1.33
1	H	96	HIS	CA-CB	5.85	1.62	1.53
2	L	59	PRO	N-CA	-5.77	1.40	1.47
1	H	208	LYS	CA-C	-5.74	1.45	1.52
2	L	92	TYR	C-O	5.69	1.30	1.24
2	L	117	ILE	N-CA	5.69	1.52	1.46
2	L	83	LEU	CA-C	5.68	1.60	1.52
1	H	107	THR	CA-C	-5.61	1.45	1.52
2	L	190	ASN	CA-C	5.60	1.56	1.52
2	L	4	MET	C-O	-5.55	1.16	1.23
2	L	58	VAL	CA-C	-5.55	1.48	1.53
1	H	124	LEU	C-O	-5.54	1.17	1.23
2	L	122	SER	CA-C	5.52	1.60	1.52
1	H	99	TYR	C-O	5.52	1.30	1.23
2	L	27	GLN	CB-CG	5.52	1.69	1.52
2	L	54	ARG	CA-C	-5.50	1.46	1.52
1	H	74	SER	CA-CB	5.47	1.62	1.53
1	H	182	THR	N-CA	5.42	1.52	1.46
2	L	87	TYR	CA-C	-5.38	1.46	1.52
1	H	27	PHE	C-O	5.34	1.30	1.23
2	L	150	ILE	CA-CB	-5.34	1.47	1.53
2	L	122	SER	C-O	-5.32	1.17	1.24
2	L	93	ASN	CG-ND2	-5.31	1.22	1.33
1	H	201	ALA	CA-C	-5.31	1.45	1.52
2	L	78	VAL	N-CA	5.27	1.52	1.46
2	L	107	LYS	CA-CB	-5.23	1.45	1.53
1	H	95	ASP	N-CA	5.21	1.52	1.46
1	H	193	VAL	CA-CB	-5.19	1.48	1.54
2	L	13	VAL	CA-C	-5.17	1.46	1.52
2	L	2	ILE	CA-CB	5.17	1.60	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	3	VAL	N-CA	5.12	1.52	1.46
2	L	138	ASN	CG-OD1	-5.10	1.13	1.23
2	L	203	SER	CA-C	5.09	1.58	1.52
1	H	196	ASN	CB-CG	5.08	1.64	1.52
1	H	208	LYS	N-CA	5.08	1.52	1.46
2	L	131	SER	C-O	-5.07	1.18	1.24
1	H	4	LEU	CG-CD2	-5.07	1.35	1.52
1	H	40	PRO	CA-C	-5.05	1.49	1.51
1	H	123	PRO	C-O	5.05	1.29	1.23
2	L	93	ASN	N-CA	5.03	1.52	1.46
2	L	206	VAL	N-CA	5.02	1.52	1.46
1	H	187	THR	CB-OG1	5.02	1.51	1.43

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	27	GLN	CA-CB-CG	9.83	133.76	114.10
1	H	19	ARG	NE-CZ-NH1	8.96	130.46	121.50
1	H	95	ASP	CA-C-N	7.90	132.04	120.90
1	H	95	ASP	C-N-CA	7.90	132.04	120.90
2	L	21	MET	CG-SD-CE	7.55	117.51	100.90
1	H	39	GLN	O-C-N	7.44	125.74	121.27
1	H	188	TRP	CB-CA-C	7.34	119.89	108.88
2	L	198	HIS	CB-CA-C	7.25	124.54	109.38
1	H	152	VAL	CA-C-N	-7.20	110.20	123.01
1	H	152	VAL	C-N-CA	-7.20	110.20	123.01
2	L	81	GLU	CB-CG-CD	7.15	124.76	112.60
1	H	208	LYS	CA-C-O	7.05	127.71	120.24
1	H	153	THR	CA-CB-OG1	-6.98	99.12	109.60
1	H	19	ARG	CD-NE-CZ	6.84	133.98	124.40
1	H	196	ASN	CA-C-O	-6.70	112.94	120.31
1	H	73	ASN	O-C-N	-6.66	112.70	122.43
1	H	159	LEU	CA-C-O	-6.64	112.94	120.50
1	H	83	ARG	NE-CZ-NH2	-6.62	113.24	119.20
2	L	48	ILE	CA-CB-CG1	6.48	121.42	110.40
2	L	198	HIS	ND1-CE1-NE2	-6.47	101.93	108.40
1	H	98	GLY	N-CA-C	-6.42	107.04	114.69
1	H	153	THR	CA-CB-CG2	6.37	121.33	110.50
2	L	200[A]	THR	O-C-N	-6.32	114.19	122.59
2	L	200[B]	THR	O-C-N	-6.32	114.19	122.59
1	H	31	ASP	CA-CB-CG	6.28	118.88	112.60
1	H	99	TYR	CA-C-N	-6.26	112.82	122.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	99	TYR	C-N-CA	-6.26	112.82	122.09
2	L	188	ARG	NE-CZ-NH1	6.21	127.71	121.50
1	H	97	ASP	CA-C-N	-6.14	113.89	122.86
1	H	97	ASP	C-N-CA	-6.14	113.89	122.86
2	L	198	HIS	CB-CG-CD2	6.14	139.19	131.20
2	L	197	THR	O-C-N	6.11	130.48	123.27
1	H	188	TRP	N-CA-C	-6.09	100.95	110.14
1	H	206	VAL	CA-C-N	-6.07	114.43	122.99
1	H	206	VAL	C-N-CA	-6.07	114.43	122.99
2	L	137	ASN	CA-CB-CG	-6.04	106.56	112.60
1	H	183	VAL	CA-CB-CG2	6.04	120.67	110.40
1	H	153	THR	O-C-N	6.02	131.05	122.74
2	L	27	GLN	CB-CG-CD	5.96	122.74	112.60
1	H	196	ASN	O-C-N	5.86	130.18	123.27
2	L	198	HIS	CB-CG-ND1	-5.83	113.95	122.70
1	H	99	TYR	O-C-N	-5.75	116.16	121.79
1	H	158[A]	SER	CA-C-N	5.64	131.42	122.74
1	H	158[A]	SER	C-N-CA	5.64	131.42	122.74
1	H	158[B]	SER	CA-C-N	5.64	131.42	122.74
1	H	158[B]	SER	C-N-CA	5.64	131.42	122.74
2	L	206	VAL	CA-CB-CG1	5.62	119.95	110.40
1	H	189	PRO	O-C-N	5.61	131.76	121.10
2	L	156	GLN	CA-C-O	5.53	125.66	119.41
1	H	118	PRO	CA-C-N	-5.52	114.57	120.03
1	H	118	PRO	C-N-CA	-5.52	114.57	120.03
2	L	179	LEU	CA-C-N	-5.45	115.31	123.00
2	L	179	LEU	C-N-CA	-5.45	115.31	123.00
1	H	107	THR	CB-CA-C	5.45	119.56	109.70
2	L	3	VAL	CB-CA-C	-5.38	102.66	110.63
2	L	108	ARG	NE-CZ-NH2	5.37	124.03	119.20
1	H	19	ARG	NE-CZ-NH2	-5.36	114.38	119.20
1	H	118	PRO	CB-CA-C	-5.30	104.45	110.92
2	L	175[A]	MET	CG-SD-CE	5.30	112.57	100.90
2	L	175[B]	MET	CG-SD-CE	5.30	112.57	100.90
1	H	194	THR	CB-CA-C	-5.21	100.27	109.70
2	L	200[A]	THR	CA-C-N	5.18	131.43	121.54
2	L	200[A]	THR	C-N-CA	5.18	131.43	121.54
2	L	200[B]	THR	CA-C-N	5.18	131.43	121.54
2	L	200[B]	THR	C-N-CA	5.18	131.43	121.54
2	L	198	HIS	N-CA-C	-5.16	101.34	109.50
1	H	95	ASP	CB-CA-C	5.15	120.66	110.42
1	H	99	TYR	CA-C-O	5.11	125.02	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	52	ARG	NE-CZ-NH2	-5.09	114.62	119.20
2	L	17	GLU	CG-CD-OE1	5.08	130.08	118.40
1	H	211	VAL	CB-CA-C	5.08	120.60	111.36
1	H	95	ASP	CA-CB-CG	5.06	117.66	112.60
1	H	39	GLN	CA-C-O	-5.05	115.33	120.63
2	L	123	GLU	N-CA-C	-5.04	105.86	111.36
2	L	127	SER	N-CA-C	5.01	119.08	112.92
2	L	181	LEU	CD1-CG-CD2	5.00	121.81	110.80
1	H	192	THR	CA-C-O	5.00	126.41	120.96

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	194	THR	Mainchain
1	H	73	ASN	Mainchain
1	H	9	GLY	Mainchain
2	L	25[A]	SER	Mainchain

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	H	1731	0	1661	34	0
2	L	1761	0	1701	18	0
3	H	3	0	0	5	0
3	L	2	0	0	0	0
4	H	319	0	0	17	0
4	L	273	0	0	10	0
All	All	4089	0	3362	54	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 8.

All (54) 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:H:169[A]:VAL:HG21	4:L:465:HOH:O	1.31	1.28
1:H:100(A):GLU:N	4:H:401:HOH:O	1.71	1.18
2:L:137:ASN:OD1	4:L:401:HOH:O	1.75	1.02
1:H:194:THR:HG23	4:H:402:HOH:O	1.67	0.93
1:H:155[A]:ASN:C	4:H:403[A]:HOH:O	2.17	0.88
2:L:81:GLU:OE2	4:L:402:HOH:O	1.92	0.87
2:L:27:GLN:HG3	4:L:524:HOH:O	1.75	0.87
1:H:100(A):GLU:C	4:H:401:HOH:O	2.24	0.81
2:L:185:GLU:OE2	2:L:188:ARG:NH2	2.16	0.79
1:H:169[A]:VAL:HG22	4:H:588:HOH:O	1.83	0.78
2:L:190:ASN:HD21	2:L:212:ASN:HB3	1.47	0.78
1:H:204:THR:HG22	4:H:516:HOH:O	1.83	0.76
1:H:156[A]:SER:N	4:H:403[A]:HOH:O	2.18	0.75
1:H:156[A]:SER:N	1:H:196:ASN:OD1	2.18	0.75
2:L:17:GLU:OE1	4:L:403:HOH:O	2.06	0.72
1:H:169[A]:VAL:CG2	4:L:465:HOH:O	2.08	0.71
2:L:13:VAL:HG13	2:L:17:GLU:HB3	1.70	0.71
3:H:302:BR:BR	4:H:621:HOH:O	2.64	0.70
2:L:27:GLN:CG	4:L:524:HOH:O	2.37	0.70
1:H:155[A]:ASN:CB	1:H:158[A]:SER:HB2	2.21	0.69
1:H:117:THR:HG22	4:H:407:HOH:O	1.94	0.67
1:H:99:TYR:CE1	4:H:629:HOH:O	2.50	0.65
1:H:194:THR:CG2	4:H:402:HOH:O	2.27	0.64
1:H:151:THR:OG1	1:H:198:ALA:HB3	1.99	0.62
1:H:155[A]:ASN:HB2	1:H:158[A]:SER:HB2	1.82	0.62
1:H:156[A]:SER:H	1:H:196:ASN:CG	2.08	0.62
1:H:157[B]:GLY:O	1:H:160:SER:OG	2.16	0.62
1:H:169[A]:VAL:CG2	4:H:588:HOH:O	2.43	0.60
1:H:155[A]:ASN:HB3	1:H:158[A]:SER:HB2	1.82	0.60
1:H:158[B]:SER:HA	4:H:410:HOH:O	2.08	0.54
2:L:13:VAL:HG11	2:L:78:VAL:HG21	1.89	0.54
1:H:19:ARG:NH2	4:H:409:HOH:O	2.41	0.53
2:L:18:LYS:HE2	4:L:622:HOH:O	2.07	0.53
3:H:302:BR:BR	4:H:484:HOH:O	2.73	0.53
1:H:100:TYR:HB2	3:H:303:BR:BR	2.64	0.52
1:H:159:LEU:O	3:H:301:BR:BR	2.83	0.52
1:H:157[B]:GLY:HA2	3:H:301:BR:BR	2.64	0.51
1:H:100(A):GLU:O	4:H:401:HOH:O	2.19	0.51
2:L:93:ASN:HD22	2:L:93:ASN:C	2.20	0.49
1:H:204:THR:CG2	4:H:516:HOH:O	2.51	0.48
1:H:188:TRP:CG	1:H:189:PRO:HA	2.51	0.46
2:L:117:ILE:HG13	2:L:134[B]:CYS:SG	2.56	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:190:ASN:HD21	2:L:212:ASN:CB	2.23	0.46
2:L:89:LYS:NZ	4:L:409:HOH:O	2.44	0.45
2:L:136:LEU:HD21	2:L:146:VAL:HG22	1.98	0.44
1:H:138:LEU:HD23	1:H:210:ILE:HG21	2.00	0.44
1:H:135:MET:HE3	1:H:182:THR:HG22	2.00	0.44
1:H:154:TRP:CZ3	1:H:195[A]:CYS:HB3	2.53	0.44
1:H:177:LEU:C	1:H:177:LEU:HD12	2.44	0.43
2:L:136:LEU:HD23	2:L:144:ILE:HD13	2.02	0.41
2:L:18:LYS:CE	4:L:622:HOH:O	2.68	0.41
1:H:188:TRP:HA	1:H:189:PRO:HA	1.76	0.41
1:H:156[A]:SER:H	1:H:196:ASN:ND2	2.17	0.41
2:L:175[A]:MET:SD	2:L:175[A]:MET:C	3.03	0.41

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	H	225/222 (101%)	216 (96%)	9 (4%)	0	100	100
2	L	227/219 (104%)	222 (98%)	5 (2%)	0	100	100
All	All	452/441 (102%)	438 (97%)	14 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	H	195/187 (104%)	187 (96%)	8 (4%)	27 2
2	L	203/194 (105%)	193 (95%)	10 (5%)	22 1
All	All	398/381 (104%)	380 (96%)	18 (4%)	26 2

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	41	PRO
1	H	50	PHE
1	H	99	TYR
1	H	134	SER
1	H	147	PRO
1	H	160	SER
1	H	191	GLU
1	H	205	LYS
2	L	11	LEU
2	L	13	VAL
2	L	123	GLU
2	L	147	LYS
2	L	175[A]	MET
2	L	175[B]	MET
2	L	181	LEU
2	L	199[A]	LYS
2	L	199[B]	LYS
2	L	212	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	39	GLN
1	H	164	HIS
2	L	38	GLN
2	L	93	ASN
2	L	157	ASN
2	L	190	ASN
2	L	212	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](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

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.

5.7 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	H	215/222 (96%)	0.34	19 (8%) 15 17	4, 14, 37, 58	12 (5%)
2	L	219/219 (100%)	0.19	2 (0%) 81 84	5, 16, 31, 45	10 (4%)
All	All	434/441 (98%)	0.27	21 (4%) 35 41	4, 15, 34, 58	22 (5%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	98	GLY	4.1
1	H	190	SER	3.7
1	H	99	TYR	3.5
1	H	189	PRO	3.4
1	H	159	LEU	3.4
2	L	126	THR	3.4
1	H	188	TRP	3.2
1	H	187	THR	2.9
1	H	213	ARG	2.9
1	H	100	TYR	2.9
1	H	158[A]	SER	2.6
1	H	195[A]	CYS	2.5
1	H	134	SER	2.5
1	H	155[A]	ASN	2.5
1	H	96	HIS	2.4
1	H	204	THR	2.4
1	H	169[A]	VAL	2.3
1	H	172[A]	SER	2.3
1	H	211	VAL	2.2
1	H	194	THR	2.2
2	L	212	ASN	2.2

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](#)

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($\text{\AA}^2$)	Q<0.9
3	BR	H	301	1/1	0.98	0.08	21,21,21,21	1
3	BR	H	303	1/1	0.99	0.17	19,19,19,19	1
3	BR	L	301	1/1	0.99	0.07	21,21,21,21	1
3	BR	L	302	1/1	0.99	0.04	25,25,25,25	1
3	BR	H	302	1/1	1.00	0.10	22,22,22,22	1

6.5 Other polymers [i](#)

There are no such residues in this entry.