



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:11 PM UTC

PDB ID : 1SSM / pdb_00001ssm
Title : Serine Acetyltransferase- Apoenzyme (truncated)
Authors : Olsen, L.R.; Huang, B.; Vetting, M.W.; Roderick, S.L.
Deposited on : 2004-03-24
Resolution : 2.15 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

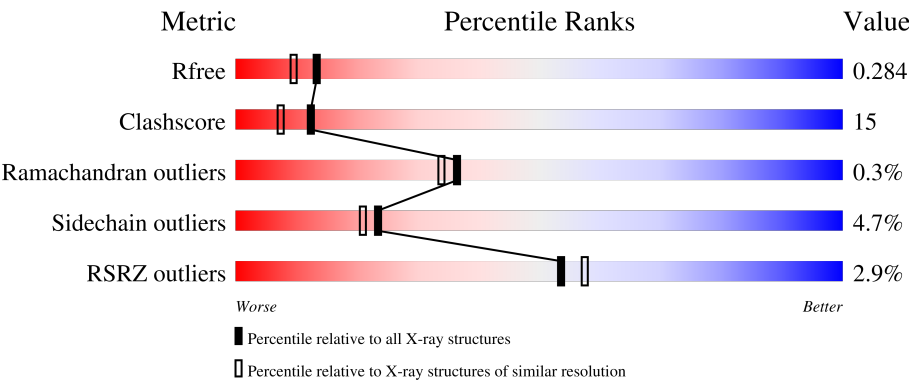
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	242	2% 64%33%..
1	B	242	3% 65%27%7%. .
1	C	242	2% 67%27%6%. .
1	D	242	3% 69%27%..
1	E	242	4% 65%29%..

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Mol	Chain	Length	Quality of chain
1	F	242	2% 60% 33% 5% **

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11514 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 Serine acetyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	S	Se	0	0	0
			1824	1162	320	336	2	4			
1	B	240	Total	C	N	O	S	Se	0	0	0
			1826	1163	319	337	2	5			
1	C	240	Total	C	N	O	S	Se	0	0	0
			1819	1159	318	336	2	4			
1	D	240	Total	C	N	O	S	Se	0	0	0
			1823	1162	319	336	2	4			
1	E	240	Total	C	N	O	S	Se	0	0	0
			1831	1166	321	337	2	5			
1	F	240	Total	C	N	O	S	Se	0	0	0
			1822	1162	319	335	2	4			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP P43886
A	22	MSE	MET	modified residue	UNP P43886
A	54	MSE	MET	modified residue	UNP P43886
A	151	MSE	MET	modified residue	UNP P43886
A	197	MSE	MET	modified residue	UNP P43886
B	1	MSE	MET	modified residue	UNP P43886
B	22	MSE	MET	modified residue	UNP P43886
B	54	MSE	MET	modified residue	UNP P43886
B	151	MSE	MET	modified residue	UNP P43886
B	197	MSE	MET	modified residue	UNP P43886
C	1	MSE	MET	modified residue	UNP P43886
C	22	MSE	MET	modified residue	UNP P43886
C	54	MSE	MET	modified residue	UNP P43886
C	151	MSE	MET	modified residue	UNP P43886
C	197	MSE	MET	modified residue	UNP P43886
D	1	MSE	MET	modified residue	UNP P43886
D	22	MSE	MET	modified residue	UNP P43886

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Chain	Residue	Modelled	Actual	Comment	Reference
D	54	MSE	MET	modified residue	UNP P43886
D	151	MSE	MET	modified residue	UNP P43886
D	197	MSE	MET	modified residue	UNP P43886
E	1	MSE	MET	modified residue	UNP P43886
E	22	MSE	MET	modified residue	UNP P43886
E	54	MSE	MET	modified residue	UNP P43886
E	151	MSE	MET	modified residue	UNP P43886
E	197	MSE	MET	modified residue	UNP P43886
F	1	MSE	MET	modified residue	UNP P43886
F	22	MSE	MET	modified residue	UNP P43886
F	54	MSE	MET	modified residue	UNP P43886
F	151	MSE	MET	modified residue	UNP P43886
F	197	MSE	MET	modified residue	UNP P43886

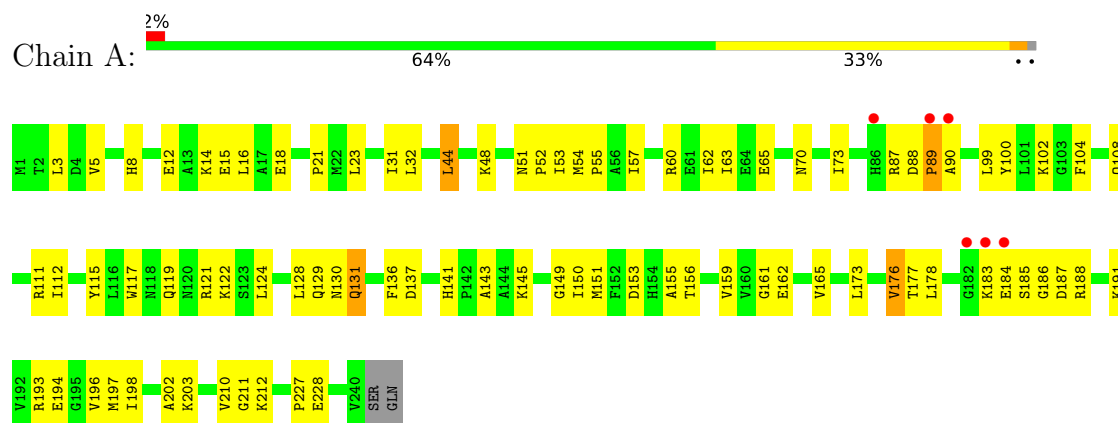
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	101	Total O 101 101	0	0
2	B	96	Total O 96 96	0	0
2	C	89	Total O 89 89	0	0
2	D	91	Total O 91 91	0	0
2	E	86	Total O 86 86	0	0
2	F	106	Total O 106 106	0	0

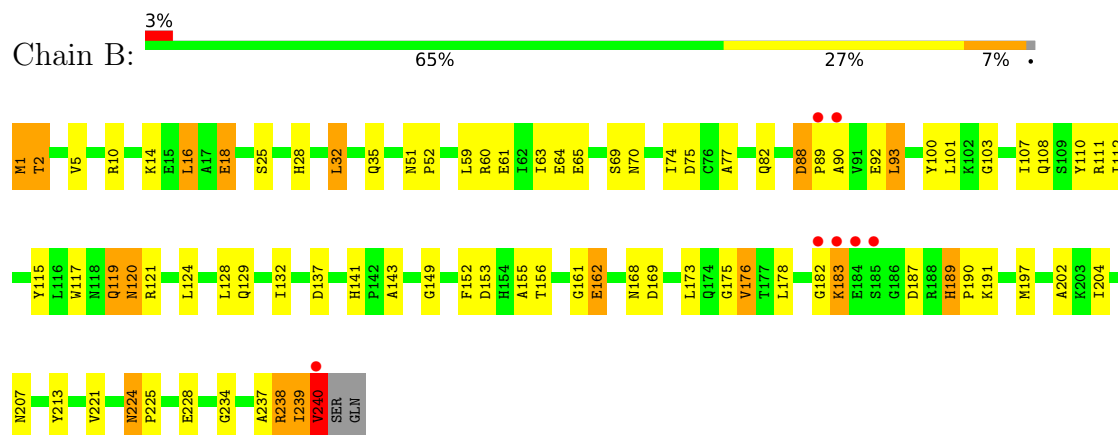
3 Residue-property plots [i](#)

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.

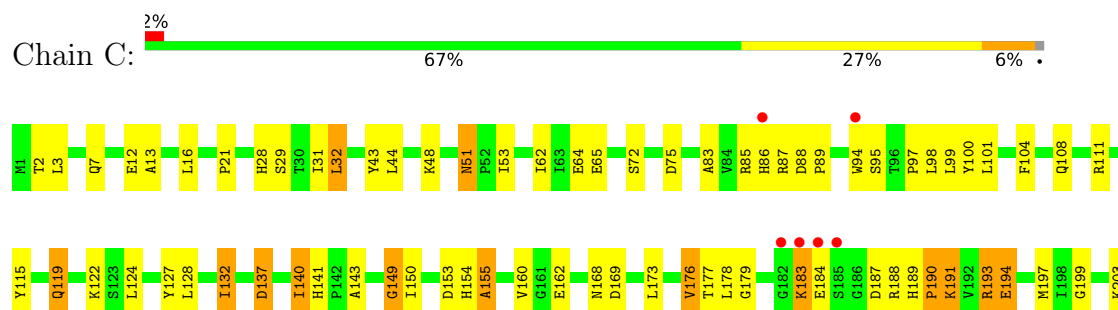
• Molecule 1: Serine acetyltransferase



• Molecule 1: Serine acetyltransferase

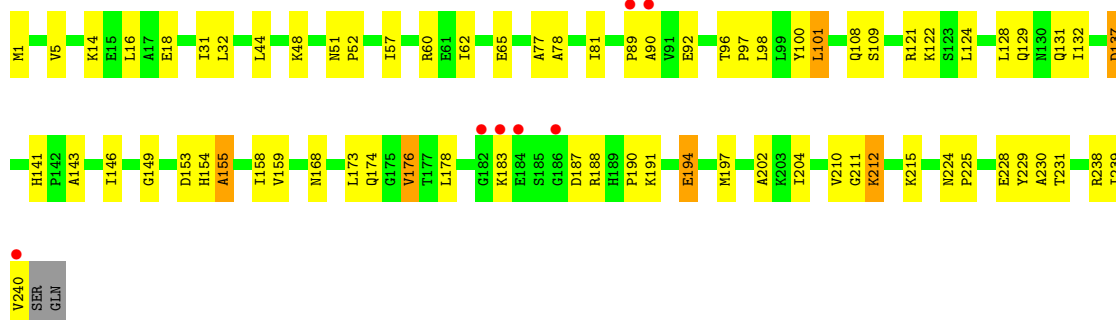


• Molecule 1: Serine acetyltransferase

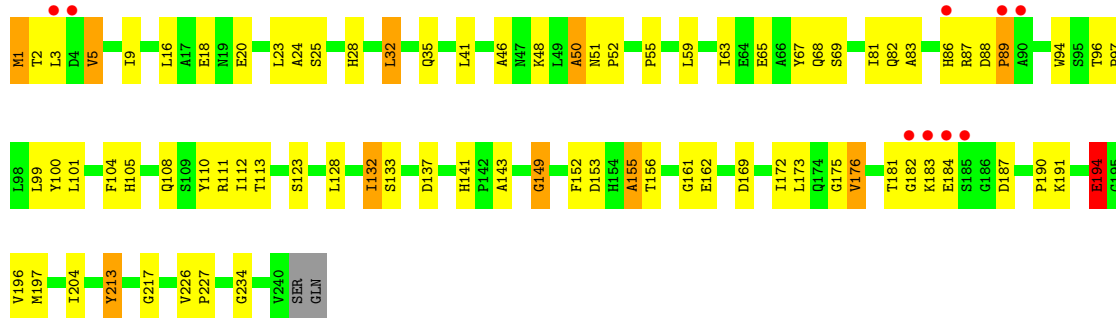




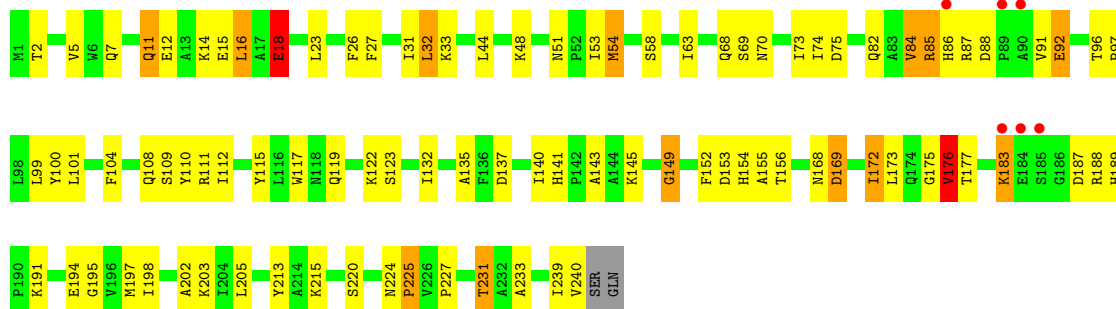
• Molecule 1: Serine acetyltransferase



• Molecule 1: Serine acetyltransferase



• Molecule 1: Serine acetyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.10Å 126.60Å 107.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.74 – 2.15 29.74 – 2.15	Depositor EDS
% Data completeness (in resolution range)	88.9 (29.74-2.15) 89.4 (29.74-2.15)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.52 (at 2.14Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.227 , 0.278 0.231 , 0.284	Depositor DCC
R_{free} test set	7181 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	22.4	Xtriage
Anisotropy	0.336	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11514	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 47.91 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.2745e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.98	0/1859	1.33	20/2525 (0.8%)
1	B	0.99	0/1860	1.40	30/2525 (1.2%)
1	C	1.03	1/1853 (0.1%)	1.31	24/2517 (1.0%)
1	D	0.98	0/1857	1.33	18/2521 (0.7%)
1	E	0.99	1/1866 (0.1%)	1.29	17/2533 (0.7%)
1	F	0.98	1/1857 (0.1%)	1.35	25/2522 (1.0%)
All	All	0.99	3/11152 (0.0%)	1.34	134/15143 (0.9%)

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	B	0	1
1	C	0	1
1	E	0	1
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	55	PRO	CA-C	6.41	1.60	1.52
1	F	176	VAL	CA-CB	5.27	1.61	1.54
1	C	88	ASP	N-CA	-5.07	1.41	1.46

All (134) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	183	LYS	N-CA-C	-12.22	91.89	110.10
1	E	5	VAL	N-CA-C	-12.12	97.79	111.00
1	B	2	THR	N-CA-C	10.62	122.43	111.07
1	F	84	VAL	N-CA-C	-9.87	100.94	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	70	ASN	CA-C-N	9.78	130.82	119.47
1	B	70	ASN	C-N-CA	9.78	130.82	119.47
1	C	32	LEU	N-CA-C	9.52	123.93	111.75
1	F	32	LEU	N-CA-C	8.94	122.15	111.33
1	F	183	LYS	N-CA-C	-8.69	96.81	109.59
1	B	93	LEU	N-CA-C	8.59	122.99	110.28
1	C	100	TYR	N-CA-C	8.34	124.14	113.88
1	D	194	GLU	N-CA-C	8.23	120.95	110.24
1	D	100	TYR	N-CA-C	8.15	123.91	113.88
1	E	161	GLY	N-CA-C	8.12	121.27	111.93
1	A	183	LYS	N-CA-C	-7.96	98.10	109.96
1	F	100	TYR	N-CA-C	7.93	123.87	114.04
1	A	131	GLN	N-CA-C	-7.89	103.10	112.89
1	A	194	GLU	N-CA-C	7.81	120.40	110.24
1	F	224	ASN	CA-C-N	7.72	127.78	119.90
1	F	224	ASN	C-N-CA	7.72	127.78	119.90
1	B	100	TYR	N-CA-C	7.45	123.28	114.04
1	D	131	GLN	N-CA-C	-7.43	103.68	112.89
1	C	194	GLU	N-CA-C	7.40	120.41	110.35
1	D	122	LYS	N-CA-C	7.21	120.19	111.82
1	E	100	TYR	N-CA-C	7.19	122.61	113.55
1	A	32	LEU	N-CA-C	7.16	121.12	112.38
1	B	155	ALA	N-CA-C	7.05	125.82	110.80
1	A	178	LEU	N-CA-C	-6.99	93.85	107.62
1	F	54	MSE	CA-C-N	6.92	126.86	120.21
1	F	54	MSE	C-N-CA	6.92	126.86	120.21
1	B	178	LEU	N-CA-C	-6.88	92.63	107.49
1	F	155	ALA	N-CA-C	6.85	125.38	110.80
1	E	32	LEU	N-CA-C	6.83	119.57	111.71
1	A	65	GLU	N-CA-C	-6.76	103.84	111.14
1	D	98	LEU	N-CA-C	6.76	119.22	111.11
1	A	149	GLY	N-CA-C	-6.75	104.69	115.08
1	D	121	ARG	N-CA-C	-6.66	100.32	109.71
1	D	101	LEU	N-CA-C	6.66	118.99	108.67
1	A	155	ALA	N-CA-C	6.62	124.91	110.80
1	B	237	ALA	N-CA-C	6.51	119.91	110.28
1	D	32	LEU	N-CA-C	6.50	120.31	112.38
1	B	169	ASP	N-CA-C	6.48	120.46	111.90
1	C	193	ARG	N-CA-C	6.43	118.16	110.19
1	B	175	GLY	N-CA-C	6.39	123.86	115.36
1	B	32	LEU	N-CA-C	6.37	119.03	111.33
1	D	211	GLY	N-CA-C	6.35	119.24	111.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	194	GLU	N-CA-C	6.31	118.45	110.24
1	F	101	LEU	N-CA-C	6.31	119.08	108.23
1	F	149	GLY	N-CA-C	-6.26	104.12	114.76
1	B	224	ASN	CA-C-N	6.24	127.64	119.84
1	B	224	ASN	C-N-CA	6.24	127.64	119.84
1	A	100	TYR	N-CA-C	6.18	122.29	114.31
1	F	169	ASP	N-CA-C	6.14	120.11	112.24
1	D	178	LEU	N-CA-C	-6.14	95.34	107.69
1	B	239	ILE	CA-C-N	6.13	132.73	121.70
1	B	239	ILE	C-N-CA	6.13	132.73	121.70
1	B	101	LEU	N-CA-C	6.12	118.15	108.67
1	E	155	ALA	N-CA-C	6.10	123.80	110.80
1	C	83	ALA	N-CA-C	6.10	117.60	111.07
1	C	183	LYS	N-CA-C	-6.08	97.85	110.80
1	B	162	GLU	N-CA-C	6.02	118.61	111.33
1	B	119	GLN	N-CA-C	-5.93	105.62	112.92
1	B	189	HIS	CA-C-N	5.91	125.92	119.89
1	B	189	HIS	C-N-CA	5.91	125.92	119.89
1	F	119	GLN	N-CA-C	-5.91	105.65	112.92
1	F	86	HIS	N-CA-C	5.87	120.30	113.20
1	D	31	ILE	N-CA-C	5.87	121.54	109.34
1	B	161	GLY	N-CA-C	5.84	120.02	112.54
1	B	121	ARG	N-CA-C	-5.79	97.76	108.24
1	A	121	ARG	N-CA-C	-5.78	101.55	109.71
1	C	168	ASN	N-CA-C	5.76	118.18	110.35
1	F	91	VAL	N-CA-C	-5.71	99.65	107.99
1	E	149	GLY	N-CA-C	-5.70	106.12	114.90
1	D	155	ALA	N-CA-C	5.69	122.93	110.80
1	C	218	ALA	N-CA-C	5.68	118.30	110.35
1	B	88	ASP	CA-C-N	5.66	125.51	119.28
1	B	88	ASP	C-N-CA	5.66	125.51	119.28
1	F	189	HIS	CA-C-N	5.65	125.60	119.78
1	F	189	HIS	C-N-CA	5.65	125.60	119.78
1	A	198	ILE	N-CA-C	5.65	115.99	107.80
1	F	205	LEU	N-CA-C	5.65	118.16	109.07
1	C	122	LYS	N-CA-C	5.64	117.51	111.36
1	A	104	PHE	N-CA-C	-5.63	105.04	111.07
1	E	99	LEU	N-CA-C	5.63	118.62	111.69
1	D	168	ASN	N-CA-C	5.61	117.57	110.33
1	B	225	PRO	CA-C-N	-5.61	118.45	123.33
1	B	225	PRO	C-N-CA	-5.61	118.45	123.33
1	C	155	ALA	N-CA-C	5.58	122.68	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	149	GLY	N-CA-C	-5.58	105.28	114.76
1	B	183	LYS	N-CA-C	-5.57	98.93	110.80
1	F	85	ARG	CA-C-N	-5.55	110.83	121.94
1	F	85	ARG	C-N-CA	-5.55	110.83	121.94
1	A	54	MSE	CA-C-N	5.53	125.84	119.92
1	A	54	MSE	C-N-CA	5.53	125.84	119.92
1	D	146	ILE	N-CA-C	5.53	115.83	107.75
1	E	101	LEU	N-CA-C	5.53	117.11	109.15
1	A	87	ARG	N-CA-C	-5.49	106.59	113.28
1	A	185	SER	N-CA-C	5.49	116.60	108.86
1	A	5	VAL	N-CA-C	-5.46	105.06	110.62
1	A	156	THR	CB-CA-C	-5.44	101.68	109.84
1	A	161	GLY	N-CA-C	5.42	119.48	112.54
1	C	101	LEU	N-CA-C	5.42	117.55	108.23
1	C	119	GLN	N-CA-C	-5.40	106.54	113.02
1	E	81	ILE	N-CA-C	-5.39	105.25	110.42
1	F	123	SER	N-CA-C	-5.38	105.41	111.28
1	E	169	ASP	N-CA-C	5.37	118.99	111.90
1	D	210	VAL	N-CA-C	-5.35	100.04	107.80
1	F	135	ALA	N-CA-C	5.31	116.88	111.14
1	F	18	GLU	CB-CA-C	-5.30	101.84	110.85
1	E	175	GLY	N-CA-C	5.28	122.91	115.30
1	E	217	GLY	N-CA-C	5.28	118.00	111.93
1	B	240	VAL	CA-C-O	-5.27	111.84	120.80
1	E	123	SER	N-CA-C	-5.24	105.74	111.82
1	C	132	ILE	N-CA-C	-5.21	105.45	110.72
1	C	190	PRO	N-CA-C	5.21	119.38	111.15
1	C	205	LEU	N-CA-C	5.20	117.38	108.90
1	C	178	LEU	N-CA-C	-5.18	97.42	107.62
1	B	149	GLY	N-CA-C	-5.17	107.11	115.08
1	C	149	GLY	N-CA-C	-5.16	105.98	114.76
1	F	5	VAL	N-CA-C	-5.16	105.47	110.42
1	E	50	ALA	N-CA-C	5.15	117.36	110.35
1	C	98	LEU	N-CA-C	5.12	116.86	111.28
1	F	175	GLY	N-CA-C	5.11	122.66	115.30
1	C	140	ILE	CB-CA-C	-5.10	104.66	110.73
1	E	234	GLY	N-CA-C	5.08	119.29	112.17
1	D	239	ILE	CB-CA-C	-5.04	105.86	111.35
1	E	132	ILE	N-CA-C	-5.03	105.64	110.72
1	A	210	VAL	N-CA-C	-5.03	99.93	107.37
1	C	65	GLU	N-CA-C	-5.03	105.90	112.23
1	B	221	VAL	N-CA-C	-5.02	100.42	107.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	51	ASN	CA-C-N	5.02	125.04	119.32
1	C	51	ASN	C-N-CA	5.02	125.04	119.32
1	C	189	HIS	CA-C-N	5.01	125.01	119.90
1	C	189	HIS	C-N-CA	5.01	125.01	119.90

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	213	TYR	Sidechain
1	C	127	TYR	Sidechain
1	E	213	TYR	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	A	1824	0	1821	53	0
1	B	1826	0	1829	72	0
1	C	1819	0	1816	56	0
1	D	1823	0	1827	54	0
1	E	1831	0	1834	56	0
1	F	1822	0	1823	67	0
2	A	101	0	0	15	0
2	B	96	0	0	14	0
2	C	89	0	0	11	0
2	D	91	0	0	9	0
2	E	86	0	0	4	0
2	F	106	0	0	9	0
All	All	11514	0	10950	338	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 15.

All (338) 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:53:ILE:HB	2:A:7567:HOH:O	1.47	1.12
1:E:141:HIS:HD2	1:E:143:ALA:H	1.00	0.98
1:E:3:LEU:HG	2:E:7410:HOH:O	1.65	0.95
1:E:141:HIS:CD2	1:E:143:ALA:H	1.85	0.94
1:C:141:HIS:HD2	1:C:143:ALA:H	1.04	0.94
1:B:141:HIS:HD2	1:B:143:ALA:H	1.00	0.93
1:D:141:HIS:HD2	1:D:143:ALA:H	1.11	0.91
1:B:92:GLU:HG3	1:B:93:LEU:H	1.37	0.90
1:B:141:HIS:CD2	1:B:143:ALA:H	1.90	0.89
1:F:197:MSE:HE2	1:F:215:LYS:HG2	1.54	0.89
1:D:141:HIS:CD2	1:D:143:ALA:H	1.91	0.88
1:C:141:HIS:CD2	1:C:143:ALA:H	1.92	0.87
1:C:3:LEU:O	1:C:7:GLN:HG2	1.75	0.86
1:B:238:ARG:NH1	1:B:238:ARG:HB3	1.93	0.83
1:B:92:GLU:HG3	1:B:93:LEU:HG	1.58	0.81
1:C:21:PRO:HG3	2:C:7555:HOH:O	1.80	0.81
1:C:31:ILE:HD11	1:C:99:LEU:HD21	1.61	0.80
1:B:92:GLU:HG3	1:B:93:LEU:N	1.95	0.80
1:F:7:GLN:NE2	2:F:7365:HOH:O	2.18	0.77
1:B:59:LEU:HD11	2:B:7465:HOH:O	1.85	0.76
1:E:41:LEU:HD23	1:E:67:TYR:OH	1.86	0.76
1:D:14:LYS:O	1:D:18:GLU:HG2	1.86	0.76
1:E:141:HIS:HD2	1:E:143:ALA:N	1.83	0.74
1:B:88:ASP:OD1	1:B:88:ASP:C	2.28	0.74
1:D:1:MSE:O	1:D:5:VAL:HG23	1.88	0.73
1:D:89:PRO:HG2	1:E:184:GLU:CB	2.19	0.73
1:F:141:HIS:HD2	1:F:143:ALA:H	1.38	0.72
1:B:187:ASP:OD1	1:B:191:LYS:NZ	2.22	0.71
1:A:184:GLU:CB	1:C:89:PRO:HG3	2.20	0.71
1:D:212:LYS:HG2	2:D:7323:HOH:O	1.89	0.71
1:D:89:PRO:HD2	2:D:7533:HOH:O	1.89	0.71
1:A:141:HIS:HD2	1:A:143:ALA:H	1.35	0.71
1:C:141:HIS:HD2	1:C:143:ALA:N	1.85	0.70
1:B:14:LYS:O	1:B:18:GLU:HG2	1.90	0.70
1:B:128:LEU:O	1:B:132:ILE:HG13	1.91	0.70
1:A:3:LEU:HD23	2:A:7544:HOH:O	1.92	0.70
1:F:197:MSE:HE1	1:F:215:LYS:NZ	2.06	0.70
1:F:231:THR:HG22	1:F:240:VAL:HG12	1.73	0.70
1:D:44:LEU:O	1:D:48:LYS:HG2	1.92	0.69
1:B:1:MSE:HG2	1:B:2:THR:N	2.08	0.69
1:A:187:ASP:OD1	1:A:191:LYS:NZ	2.25	0.69
1:D:197:MSE:HE1	1:D:215:LYS:HE2	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:153:ASP:HB3	1:F:173:LEU:HD23	1.76	0.68
1:A:141:HIS:CD2	1:A:143:ALA:H	2.10	0.68
1:E:28:HIS:NE2	2:E:7230:HOH:O	2.25	0.68
1:E:88:ASP:OD1	1:E:88:ASP:C	2.35	0.67
1:F:74:ILE:HD13	2:F:7261:HOH:O	1.93	0.67
1:A:14:LYS:O	1:A:18:GLU:HG2	1.94	0.67
1:D:173:LEU:O	1:D:176:VAL:HG13	1.95	0.66
1:E:1:MSE:HE3	1:E:2:THR:H	1.61	0.66
1:B:92:GLU:CG	1:B:93:LEU:H	2.07	0.66
1:A:90:ALA:N	2:A:7552:HOH:O	2.24	0.65
1:A:44:LEU:O	1:A:48:LYS:HG2	1.95	0.65
1:B:168:ASN:OD1	2:B:7374:HOH:O	2.14	0.65
1:C:85:ARG:NH1	1:C:94:TRP:CZ2	2.65	0.65
1:E:89:PRO:HA	2:E:7549:HOH:O	1.97	0.65
1:C:183:LYS:CB	2:C:7476:HOH:O	2.44	0.64
1:D:155:ALA:O	1:D:158:ILE:HG13	1.97	0.64
1:A:89:PRO:HD2	2:A:7552:HOH:O	1.97	0.64
1:C:31:ILE:CD1	1:C:99:LEU:HD21	2.27	0.64
1:F:141:HIS:CD2	1:F:143:ALA:H	2.15	0.64
1:B:137:ASP:OD1	1:B:156:THR:HG23	1.97	0.63
1:F:122:LYS:HG3	2:F:7300:HOH:O	1.97	0.63
1:D:96:THR:HB	1:D:97:PRO:HD3	1.79	0.63
1:B:1:MSE:HG2	1:B:2:THR:H	1.62	0.63
1:E:108:GLN:HE22	1:E:111:ARG:HE	1.47	0.63
1:A:62:ILE:HD13	1:A:124:LEU:HD11	1.81	0.63
1:F:197:MSE:HE1	1:F:215:LYS:HZ2	1.62	0.62
1:B:156:THR:HG23	1:C:137:ASP:OD2	2.00	0.62
1:C:191:LYS:HG2	1:C:193:ARG:NH2	2.14	0.62
1:A:153:ASP:HB3	1:A:173:LEU:HD23	1.81	0.62
1:C:173:LEU:O	1:C:176:VAL:HG13	2.01	0.61
1:B:108:GLN:HE22	1:B:111:ARG:HE	1.48	0.61
1:B:238:ARG:HB3	1:B:238:ARG:HH11	1.64	0.61
1:F:88:ASP:C	1:F:88:ASP:OD1	2.43	0.60
1:D:141:HIS:HD2	1:D:143:ALA:N	1.92	0.60
1:C:173:LEU:HD12	1:C:199:GLY:HA2	1.81	0.60
1:E:82:GLN:CG	1:E:94:TRP:HZ3	2.14	0.60
1:B:238:ARG:HD2	2:B:7540:HOH:O	2.02	0.60
1:B:240:VAL:HG23	1:B:240:VAL:O	2.00	0.60
1:B:108:GLN:HA	1:B:108:GLN:NE2	2.17	0.60
1:E:197:MSE:HG3	2:E:7493:HOH:O	2.02	0.59
1:F:197:MSE:HE2	1:F:215:LYS:CG	2.31	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:PRO:HG3	1:C:184:GLU:O	2.01	0.59
1:B:141:HIS:HD2	1:B:143:ALA:N	1.85	0.59
1:F:168:ASN:OD1	2:F:7332:HOH:O	2.16	0.59
1:C:115:TYR:O	1:C:119:GLN:HG2	2.03	0.59
1:C:128:LEU:O	1:C:132:ILE:HG13	2.01	0.59
1:A:159:VAL:HG11	1:C:154:HIS:CD2	2.38	0.59
1:F:122:LYS:HE3	2:F:7417:HOH:O	2.03	0.59
1:C:86:HIS:C	2:C:7513:HOH:O	2.46	0.59
1:F:183:LYS:O	1:F:188:ARG:HD2	2.03	0.59
1:B:65:GLU:OE1	1:B:69:SER:OG	2.21	0.58
1:F:70:ASN:HB3	1:F:73:ILE:HD12	1.84	0.58
1:A:14:LYS:HE3	2:A:7479:HOH:O	2.04	0.58
1:F:197:MSE:CE	1:F:215:LYS:HG2	2.31	0.58
1:A:51:ASN:ND2	2:A:7165:HOH:O	2.36	0.58
1:B:153:ASP:HB3	1:B:173:LEU:HD23	1.86	0.58
1:B:238:ARG:CD	2:B:7489:HOH:O	2.52	0.57
1:B:88:ASP:OD1	1:B:90:ALA:N	2.37	0.57
1:D:65:GLU:HB3	2:D:7346:HOH:O	2.03	0.57
1:F:44:LEU:O	1:F:48:LYS:HG2	2.03	0.57
1:B:190:PRO:CG	1:B:204:ILE:HG22	2.34	0.57
1:A:70:ASN:ND2	1:A:73:ILE:HG13	2.19	0.57
1:F:23:LEU:O	1:F:26:PHE:HB3	2.05	0.57
1:F:109:SER:HB3	1:F:132:ILE:HD11	1.87	0.57
1:C:183:LYS:O	1:C:188:ARG:HD2	2.05	0.57
1:F:187:ASP:OD1	1:F:191:LYS:NZ	2.29	0.56
1:B:92:GLU:HG2	2:B:7009:HOH:O	2.04	0.56
1:E:82:GLN:HG2	1:E:94:TRP:HZ3	1.70	0.56
1:E:156:THR:CG2	1:F:137:ASP:OD2	2.53	0.56
1:B:137:ASP:OD1	1:C:137:ASP:OD2	2.23	0.56
1:C:7:GLN:HA	1:C:7:GLN:NE2	2.20	0.56
1:D:231:THR:HG22	1:D:240:VAL:CG2	2.36	0.56
1:D:137:ASP:CG	1:F:137:ASP:OD1	2.49	0.56
1:F:183:LYS:N	2:F:7473:HOH:O	2.37	0.56
1:E:187:ASP:OD1	1:E:191:LYS:NZ	2.33	0.56
1:D:228:GLU:HG2	2:D:7338:HOH:O	2.05	0.56
1:B:238:ARG:HG2	2:B:7540:HOH:O	2.06	0.56
1:E:82:GLN:HG2	1:E:94:TRP:CZ3	2.40	0.56
1:F:173:LEU:O	1:F:176:VAL:HG13	2.06	0.56
1:A:23:LEU:HD22	1:B:124:LEU:HD12	1.86	0.55
1:B:60:ARG:O	1:B:64:GLU:HG3	2.06	0.55
1:B:28:HIS:HA	1:B:32:LEU:HB2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:GLN:HE22	1:C:111:ARG:HE	1.55	0.55
1:B:74:ILE:O	1:B:77:ALA:HB3	2.05	0.55
1:B:238:ARG:HD3	2:B:7489:HOH:O	2.07	0.55
1:F:12:GLU:O	1:F:16:LEU:HD22	2.07	0.55
1:F:177:THR:HB	1:F:203:LYS:HD2	1.88	0.54
1:B:238:ARG:HB3	1:B:238:ARG:CZ	2.36	0.54
1:D:154:HIS:O	1:D:155:ALA:HB3	2.07	0.54
1:F:225:PRO:HG3	2:F:7127:HOH:O	2.07	0.54
1:D:228:GLU:CG	2:D:7338:HOH:O	2.56	0.54
1:D:51:ASN:ND2	2:D:7215:HOH:O	2.41	0.54
1:A:108:GLN:HA	1:A:108:GLN:NE2	2.23	0.53
1:D:197:MSE:HE3	1:D:215:LYS:HG2	1.90	0.53
1:E:156:THR:HG23	1:F:137:ASP:OD2	2.08	0.53
1:E:190:PRO:CG	1:E:204:ILE:HG22	2.38	0.53
1:A:115:TYR:O	1:A:119:GLN:HG2	2.08	0.53
1:D:137:ASP:OD1	1:F:137:ASP:OD1	2.26	0.53
1:E:28:HIS:HA	1:E:32:LEU:HB2	1.89	0.53
1:F:197:MSE:CE	1:F:215:LYS:NZ	2.72	0.53
1:A:143:ALA:HB3	1:A:162:GLU:HG3	1.90	0.52
1:A:15:GLU:HG2	2:A:7506:HOH:O	2.09	0.52
1:C:94:TRP:O	1:C:97:PRO:HD2	2.09	0.52
1:B:238:ARG:CD	2:B:7540:HOH:O	2.56	0.52
1:C:51:ASN:OD1	1:C:53:ILE:N	2.43	0.52
1:B:162:GLU:N	1:B:189:HIS:CE1	2.77	0.52
1:D:89:PRO:CD	2:D:7533:HOH:O	2.54	0.52
1:E:88:ASP:OD1	1:E:88:ASP:O	2.27	0.52
1:F:231:THR:HG22	1:F:240:VAL:CG1	2.39	0.52
1:D:190:PRO:C	1:D:191:LYS:HD2	2.35	0.52
1:E:133:SER:O	1:E:137:ASP:HA	2.10	0.52
1:B:238:ARG:CG	2:B:7540:HOH:O	2.57	0.51
1:A:173:LEU:O	1:A:176:VAL:HG13	2.10	0.51
1:C:149:GLY:O	1:C:169:ASP:HA	2.10	0.51
1:E:194:GLU:OE2	1:E:213:TYR:OH	2.19	0.51
1:C:64:GLU:HG2	2:C:7486:HOH:O	2.11	0.51
1:D:137:ASP:OD2	1:F:137:ASP:OD1	2.28	0.51
1:A:88:ASP:O	1:A:90:ALA:N	2.43	0.51
1:F:31:ILE:CD1	1:F:99:LEU:HD21	2.41	0.51
1:F:108:GLN:HE22	1:F:111:ARG:HE	1.58	0.51
1:A:8:HIS:O	1:A:12:GLU:HG3	2.11	0.50
1:F:16:LEU:HB3	1:F:27:PHE:CE1	2.46	0.50
1:C:21:PRO:CG	2:C:7555:HOH:O	2.48	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MSE:O	1:B:5:VAL:HG23	2.12	0.50
1:C:87:ARG:NH1	1:C:87:ARG:HG3	2.27	0.50
1:B:128:LEU:O	1:B:129:GLN:C	2.54	0.49
1:D:109:SER:HB3	1:D:132:ILE:HD11	1.93	0.49
1:B:2:THR:OG1	1:B:75:ASP:OD1	2.24	0.49
1:D:137:ASP:OD1	1:F:156:THR:HG23	2.12	0.49
1:D:62:ILE:HD13	1:D:124:LEU:HD11	1.95	0.49
1:B:16:LEU:HD21	1:B:93:LEU:HD13	1.94	0.49
1:B:224:ASN:ND2	2:B:7387:HOH:O	2.40	0.49
1:F:109:SER:HB3	1:F:132:ILE:CD1	2.42	0.49
1:A:193:ARG:O	1:A:196:VAL:HG23	2.12	0.49
1:B:173:LEU:O	1:B:176:VAL:HG13	2.13	0.49
1:C:153:ASP:C	1:C:155:ALA:H	2.21	0.49
1:F:51:ASN:OD1	1:F:53:ILE:N	2.45	0.49
1:A:108:GLN:HE22	1:A:111:ARG:HE	1.60	0.49
1:A:197:MSE:HG3	2:A:7112:HOH:O	2.13	0.48
1:C:13:ALA:HB2	1:C:31:ILE:HG21	1.94	0.48
1:C:177:THR:HB	1:C:203:LYS:HD2	1.94	0.48
1:D:108:GLN:HA	1:D:108:GLN:NE2	2.28	0.48
1:B:61:GLU:OE1	2:B:7384:HOH:O	2.20	0.48
1:B:197:MSE:HG3	2:B:7414:HOH:O	2.12	0.48
1:C:21:PRO:CD	2:C:7555:HOH:O	2.61	0.48
1:A:186:GLY:O	1:A:188:ARG:HG2	2.13	0.48
1:C:62:ILE:HD13	1:C:124:LEU:HD11	1.95	0.48
1:C:51:ASN:OD1	1:C:51:ASN:C	2.57	0.48
1:A:63:ILE:HG23	1:A:112:ILE:CD1	2.44	0.48
1:D:197:MSE:CE	1:D:215:LYS:HE2	2.43	0.48
1:A:57:ILE:HD11	1:D:60:ARG:HD2	1.95	0.47
1:A:228:GLU:CG	2:A:7018:HOH:O	2.62	0.47
1:A:89:PRO:CD	2:A:7552:HOH:O	2.57	0.47
1:D:51:ASN:HB2	1:D:52:PRO:CD	2.44	0.47
1:D:153:ASP:O	1:D:154:HIS:HB2	2.15	0.47
1:E:51:ASN:HB2	1:E:52:PRO:CD	2.45	0.47
1:E:128:LEU:O	1:E:132:ILE:HG13	2.14	0.47
1:F:195:GLY:HA3	1:F:213:TYR:CE2	2.49	0.47
1:F:227:PRO:HD2	1:F:239:ILE:HD12	1.95	0.47
1:C:183:LYS:HA	2:C:7568:HOH:O	2.14	0.47
1:A:211:GLY:O	1:A:212:LYS:C	2.57	0.47
1:B:25:SER:HA	1:E:25:SER:HA	1.96	0.47
1:A:130:ASN:HB3	2:C:7163:HOH:O	2.13	0.47
1:B:190:PRO:HG3	1:B:204:ILE:HG22	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:LYS:HE3	1:A:136:PHE:O	2.15	0.47
1:A:176:VAL:HA	1:A:202:ALA:O	2.15	0.47
1:D:197:MSE:HE1	1:D:215:LYS:CE	2.44	0.47
1:D:159:VAL:HG11	1:F:154:HIS:CD2	2.50	0.47
1:F:227:PRO:HD2	1:F:239:ILE:CD1	2.45	0.47
1:E:111:ARG:HG2	1:E:111:ARG:HH11	1.80	0.46
1:F:96:THR:HB	1:F:97:PRO:HD3	1.96	0.46
1:F:108:GLN:HA	1:F:108:GLN:NE2	2.30	0.46
1:A:51:ASN:HB2	1:A:52:PRO:CD	2.45	0.46
1:C:95:SER:O	1:C:99:LEU:HG	2.15	0.46
1:E:65:GLU:CD	1:E:69:SER:HG	2.21	0.46
1:E:96:THR:HB	1:E:97:PRO:HD3	1.97	0.46
1:B:92:GLU:CG	1:B:93:LEU:N	2.66	0.46
1:D:228:GLU:C	1:D:230:ALA:H	2.24	0.46
1:E:82:GLN:HG3	1:E:94:TRP:HZ3	1.80	0.46
1:A:63:ILE:HG23	1:A:112:ILE:HD11	1.97	0.46
1:E:5:VAL:O	1:E:9:ILE:HG13	2.16	0.46
1:F:11:GLN:O	1:F:15:GLU:HG2	2.15	0.46
1:B:176:VAL:HA	1:B:202:ALA:O	2.15	0.46
1:E:32:LEU:HD23	1:E:32:LEU:HA	1.78	0.46
1:E:59:LEU:O	1:E:63:ILE:HG13	2.16	0.46
1:D:96:THR:O	1:D:97:PRO:C	2.58	0.46
1:E:24:ALA:O	1:E:25:SER:C	2.58	0.46
1:F:33:LYS:NZ	2:F:7140:HOH:O	2.27	0.46
1:D:176:VAL:HA	1:D:202:ALA:O	2.16	0.45
1:A:150:ILE:HG22	1:A:151:MSE:N	2.31	0.45
1:D:137:ASP:OD1	1:F:156:THR:CG2	2.64	0.45
1:F:140:ILE:HD11	1:F:152:PHE:HE1	1.81	0.45
1:B:115:TYR:O	1:B:119:GLN:HG2	2.16	0.45
1:A:89:PRO:N	2:A:7552:HOH:O	2.49	0.45
1:C:150:ILE:O	2:C:7310:HOH:O	2.21	0.45
1:A:15:GLU:CG	2:A:7400:HOH:O	2.64	0.45
1:B:107:ILE:HD11	1:B:152:PHE:CG	2.51	0.45
1:C:153:ASP:C	1:C:155:ALA:N	2.75	0.45
1:F:92:GLU:H	1:F:92:GLU:HG2	1.55	0.45
1:F:220:SER:OG	1:F:233:ALA:HA	2.17	0.45
1:B:10:ARG:O	1:B:14:LYS:HG3	2.17	0.45
1:B:103:GLY:O	1:B:107:ILE:HG13	2.17	0.45
1:C:28:HIS:HA	1:C:32:LEU:HB2	1.99	0.45
1:C:2:THR:OG1	1:C:75:ASP:OD1	2.31	0.45
1:D:77:ALA:O	1:D:81:ILE:HG13	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:48:LYS:HE3	1:E:105:HIS:CE1	2.52	0.45
1:F:87:ARG:NH1	1:F:87:ARG:HG3	2.32	0.44
1:F:117:TRP:CH2	1:F:122:LYS:HD3	2.52	0.44
1:B:239:ILE:HG22	1:B:240:VAL:HG13	1.99	0.44
1:E:173:LEU:O	1:E:176:VAL:HG13	2.17	0.44
1:F:14:LYS:O	1:F:18:GLU:HG2	2.17	0.44
1:F:104:PHE:O	1:F:108:GLN:HG2	2.18	0.44
1:A:60:ARG:HD2	1:D:57:ILE:HD11	1.99	0.44
1:C:29:SER:HB3	1:C:43:TYR:OH	2.17	0.44
1:E:153:ASP:HB3	1:E:173:LEU:HD23	1.98	0.44
1:A:177:THR:HB	1:A:203:LYS:HD2	2.00	0.44
1:D:231:THR:HG22	1:D:240:VAL:HG23	1.98	0.44
1:E:63:ILE:HG23	1:E:112:ILE:CD1	2.47	0.44
1:C:44:LEU:O	1:C:48:LYS:HG3	2.17	0.44
1:C:140:ILE:HG23	1:C:160:VAL:HB	2.00	0.44
1:B:92:GLU:HG2	1:B:92:GLU:H	1.61	0.44
1:A:53:ILE:C	1:A:55:PRO:HD3	2.43	0.43
1:E:87:ARG:C	1:E:89:PRO:HD3	2.44	0.43
1:E:196:VAL:CG1	1:E:197:MSE:N	2.81	0.43
1:A:165:VAL:HG23	1:A:191:LYS:HE3	2.00	0.43
1:A:153:ASP:HB3	1:A:173:LEU:CD2	2.48	0.43
1:F:31:ILE:HG22	1:F:32:LEU:N	2.33	0.43
1:B:108:GLN:NE2	1:B:108:GLN:CA	2.82	0.43
1:C:31:ILE:HG22	1:C:32:LEU:N	2.32	0.43
1:C:104:PHE:O	1:C:108:GLN:HG2	2.17	0.43
1:D:187:ASP:OD1	1:D:191:LYS:NZ	2.49	0.43
1:E:83:ALA:HB2	1:E:149:GLY:HA2	2.00	0.43
1:E:137:ASP:OD1	1:F:137:ASP:OD1	2.37	0.43
1:F:2:THR:OG1	1:F:75:ASP:OD1	2.29	0.43
1:E:143:ALA:HB3	1:E:162:GLU:HG3	2.00	0.43
1:A:117:TRP:CH2	1:A:122:LYS:HD3	2.54	0.43
1:B:156:THR:CG2	1:C:137:ASP:OD2	2.65	0.43
1:E:20:GLU:HG3	1:E:23:LEU:HB2	2.00	0.43
1:D:90:ALA:N	2:D:7533:HOH:O	2.52	0.43
1:B:51:ASN:HB2	1:B:52:PRO:HD3	2.01	0.43
1:E:88:ASP:O	1:E:88:ASP:CG	2.60	0.43
1:A:53:ILE:CD1	2:A:7567:HOH:O	2.67	0.42
1:C:3:LEU:N	1:C:3:LEU:HD23	2.34	0.42
1:E:1:MSE:HE3	1:E:1:MSE:HB3	1.71	0.42
1:E:152:PHE:CE2	1:E:172:ILE:HD12	2.53	0.42
1:E:182:GLY:O	1:E:184:GLU:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:176:VAL:HA	1:F:202:ALA:O	2.19	0.42
1:A:31:ILE:HD11	1:A:99:LEU:HD21	2.00	0.42
1:B:108:GLN:HA	1:B:108:GLN:HE21	1.83	0.42
1:F:172:ILE:HD13	1:F:198:ILE:HD12	2.02	0.42
1:B:92:GLU:CG	1:B:93:LEU:HG	2.41	0.42
1:C:212:LYS:HG2	1:C:213:TYR:HD1	1.84	0.42
1:E:104:PHE:O	1:E:108:GLN:HG2	2.20	0.41
1:F:23:LEU:HD12	1:F:23:LEU:HA	1.88	0.41
1:A:128:LEU:O	1:A:129:GLN:C	2.63	0.41
1:B:51:ASN:HB2	1:B:52:PRO:CD	2.50	0.41
1:B:117:TRP:O	1:B:120:ASN:N	2.49	0.41
1:B:182:GLY:O	1:B:183:LYS:C	2.63	0.41
1:D:5:VAL:HG21	1:D:78:ALA:HB1	2.02	0.41
1:D:190:PRO:CG	1:D:204:ILE:HG22	2.50	0.41
1:F:145:LYS:HD2	2:F:7162:HOH:O	2.20	0.41
1:B:173:LEU:HD23	1:B:173:LEU:HA	1.86	0.41
1:C:197:MSE:HE3	2:C:7123:HOH:O	2.20	0.41
1:D:153:ASP:HB3	1:D:173:LEU:HD23	2.03	0.41
1:D:173:LEU:HD22	1:D:174:GLN:OE1	2.20	0.41
1:E:113:THR:HG22	1:E:128:LEU:HB2	2.01	0.41
1:F:54:MSE:HE3	1:F:54:MSE:HB3	1.94	0.41
1:A:196:VAL:HG12	1:A:197:MSE:N	2.36	0.41
1:A:228:GLU:CD	2:A:7018:HOH:O	2.62	0.41
1:C:187:ASP:OD1	1:C:191:LYS:NZ	2.52	0.41
1:E:137:ASP:OD1	1:E:156:THR:HG23	2.20	0.41
1:B:234:GLY:HA2	2:B:7497:HOH:O	2.21	0.41
1:E:153:ASP:C	1:E:155:ALA:N	2.79	0.41
1:B:228:GLU:HG3	2:B:7398:HOH:O	2.19	0.41
1:C:12:GLU:OE1	1:C:94:TRP:HD1	2.03	0.41
1:D:224:ASN:HA	1:D:225:PRO:HD3	1.95	0.41
1:B:63:ILE:HG23	1:B:112:ILE:CD1	2.51	0.41
1:D:1:MSE:CB	2:D:7468:HOH:O	2.69	0.41
1:D:194:GLU:O	1:D:212:LYS:HA	2.21	0.41
1:C:141:HIS:HB3	1:C:162:GLU:HA	2.02	0.41
1:C:179:GLY:O	1:C:190:PRO:HD3	2.21	0.41
1:F:149:GLY:O	1:F:169:ASP:HA	2.20	0.41
1:B:207:ASN:O	1:B:207:ASN:CG	2.64	0.40
1:E:46:ALA:O	1:E:50:ALA:HB2	2.21	0.40
1:E:226:VAL:HA	1:E:227:PRO:HD3	1.81	0.40
1:D:97:PRO:HA	1:D:101:LEU:HG	2.02	0.40
1:D:128:LEU:O	1:D:129:GLN:C	2.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:155:ALA:O	1:E:156:THR:C	2.64	0.40
1:F:63:ILE:HG23	1:F:112:ILE:CD1	2.52	0.40
1:A:145:LYS:HG3	2:A:7345:HOH:O	2.21	0.40
1:C:32:LEU:HA	1:C:32:LEU:HD23	1.86	0.40
1:E:153:ASP:O	1:E:155:ALA:N	2.48	0.40
1:F:84:VAL:O	1:F:85:ARG:C	2.63	0.40
1:C:184:GLU:N	2:C:7568:HOH:O	2.38	0.40
1:D:191:LYS:HD2	1:D:191:LYS:N	2.36	0.40
1:F:69:SER:HB2	1:F:115:TYR:CE1	2.56	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	
1	A	238/242 (98%)	229 (96%)	8 (3%)	1 (0%)	30	26
1	B	238/242 (98%)	225 (94%)	13 (6%)	0	100	100
1	C	238/242 (98%)	229 (96%)	9 (4%)	0	100	100
1	D	238/242 (98%)	224 (94%)	13 (6%)	1 (0%)	30	26
1	E	238/242 (98%)	225 (94%)	11 (5%)	2 (1%)	16	10
1	F	238/242 (98%)	227 (95%)	11 (5%)	0	100	100
All	All	1428/1452 (98%)	1359 (95%)	65 (5%)	4 (0%)	36	34

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	229	TYR
1	E	183	LYS
1	E	89	PRO
1	A	89	PRO

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	190/192 (99%)	183 (96%)	7 (4%)	30	30
1	B	191/192 (100%)	181 (95%)	10 (5%)	21	17
1	C	189/192 (98%)	182 (96%)	7 (4%)	30	30
1	D	190/192 (99%)	183 (96%)	7 (4%)	30	30
1	E	192/192 (100%)	182 (95%)	10 (5%)	21	17
1	F	190/192 (99%)	177 (93%)	13 (7%)	14	10
All	All	1142/1152 (99%)	1088 (95%)	54 (5%)	23	21

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	21	PRO
1	A	44	LEU
1	A	131	GLN
1	A	137	ASP
1	A	176	VAL
1	A	227	PRO
1	B	1	MSE
1	B	16	LEU
1	B	18	GLU
1	B	35	GLN
1	B	82	GLN
1	B	110	TYR
1	B	120	ASN
1	B	176	VAL
1	B	238	ARG
1	B	240	VAL
1	C	16	LEU
1	C	72	SER
1	C	137	ASP
1	C	176	VAL
1	C	191	LYS

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Mol	Chain	Res	Type
1	C	194	GLU
1	C	238	ARG
1	D	16	LEU
1	D	92	GLU
1	D	137	ASP
1	D	176	VAL
1	D	188	ARG
1	D	212	LYS
1	D	238	ARG
1	E	1	MSE
1	E	16	LEU
1	E	18	GLU
1	E	35	GLN
1	E	68	GLN
1	E	86	HIS
1	E	110	TYR
1	E	176	VAL
1	E	181	THR
1	E	194	GLU
1	F	11	GLN
1	F	16	LEU
1	F	18	GLU
1	F	58	SER
1	F	68	GLN
1	F	82	GLN
1	F	92	GLU
1	F	110	TYR
1	F	172	ILE
1	F	176	VAL
1	F	194	GLU
1	F	225	PRO
1	F	231	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	68	GLN
1	A	108	GLN
1	A	130	ASN
1	A	141	HIS
1	B	7	GLN

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Mol	Chain	Res	Type
1	B	82	GLN
1	B	108	GLN
1	B	130	ASN
1	B	141	HIS
1	B	154	HIS
1	C	7	GLN
1	C	108	GLN
1	C	130	ASN
1	C	141	HIS
1	C	148	HIS
1	C	154	HIS
1	C	224	ASN
1	D	68	GLN
1	D	130	ASN
1	D	141	HIS
1	D	224	ASN
1	E	11	GLN
1	E	19	ASN
1	E	35	GLN
1	E	108	GLN
1	E	141	HIS
1	E	207	ASN
1	F	7	GLN
1	F	82	GLN
1	F	86	HIS
1	F	108	GLN
1	F	129	GLN
1	F	141	HIS
1	F	154	HIS
1	F	207	ASN
1	F	224	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	235/242 (97%)	0.10	6 (2%) 57 61	16, 25, 41, 56	0
1	B	235/242 (97%)	0.15	7 (2%) 52 56	16, 25, 41, 55	0
1	C	235/242 (97%)	0.19	6 (2%) 57 61	17, 25, 42, 52	0
1	D	235/242 (97%)	0.10	7 (2%) 52 56	15, 24, 40, 50	0
1	E	235/242 (97%)	0.16	9 (3%) 44 49	16, 25, 42, 52	0
1	F	235/242 (97%)	0.08	6 (2%) 57 61	16, 23, 41, 53	0
All	All	1410/1452 (97%)	0.13	41 (2%) 53 57	15, 25, 42, 56	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	183	LYS	6.1
1	C	184	GLU	4.8
1	E	184	GLU	4.7
1	B	184	GLU	4.7
1	C	182	GLY	4.6
1	A	184	GLU	4.3
1	B	240	VAL	4.2
1	A	183	LYS	4.1
1	A	90	ALA	3.8
1	F	86	HIS	3.8
1	D	89	PRO	3.6
1	D	183	LYS	3.5
1	A	89	PRO	3.3
1	B	89	PRO	3.3
1	F	90	ALA	3.3
1	E	4	ASP	3.2
1	D	186	GLY	3.2
1	B	182	GLY	3.1
1	E	182	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	185	SER	3.0
1	B	183	LYS	2.9
1	E	183	LYS	2.9
1	F	183	LYS	2.9
1	B	90	ALA	2.9
1	A	86	HIS	2.9
1	E	90	ALA	2.8
1	D	240	VAL	2.7
1	A	182	GLY	2.6
1	D	184	GLU	2.6
1	E	185	SER	2.5
1	F	89	PRO	2.4
1	D	90	ALA	2.3
1	C	185	SER	2.3
1	E	86	HIS	2.3
1	C	94	TRP	2.3
1	D	182	GLY	2.2
1	F	185	SER	2.2
1	E	89	PRO	2.1
1	F	184	GLU	2.1
1	E	3	LEU	2.1
1	C	86	HIS	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.