



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 04:58 PM UTC

PDB ID : 4XS6 / pdb_00004xs6
Title : Salmonella typhimurium AhpC W81F mutant
Authors : Perkins, A.; Nelson, K.; Parsonage, D.; Poole, L.; Karplus, P.A.
Deposited on : 2015-01-21
Resolution : 3.35 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

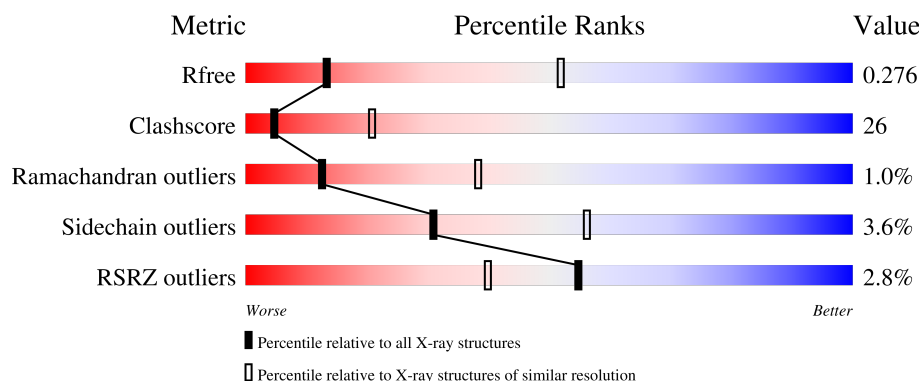
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.35 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	1099 (3.40-3.32)
Clashscore	190562	1116 (3.40-3.32)
Ramachandran outliers	187476	1101 (3.40-3.32)
Sidechain outliers	187428	1101 (3.40-3.32)
RSRZ outliers	180081	1099 (3.40-3.32)

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	186	2% 59% 24% • • 12%
1	B	186	2% 59% 25% • • 12%
1	C	186	0% 67% 20% • 13%
1	D	186	0% 57% 26% • 13%
1	E	186	7% 39% 30% 12% 7% 13%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6410 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 Alkyl hydroperoxide reductase subunit C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	163	Total	C	N	O	S	0	1	0
			1291	820	215	253	3			
1	B	163	Total	C	N	O	S	0	1	0
			1285	817	211	254	3			
1	C	162	Total	C	N	O	S	0	0	0
			1274	811	212	248	3			
1	D	162	Total	C	N	O	S	0	1	0
			1282	815	214	250	3			
1	E	162	Total	C	N	O	S	0	0	0
			1274	811	212	248	3			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	81	PHE	TRP	engineered mutation	UNP P0A251
B	81	PHE	TRP	engineered mutation	UNP P0A251
C	81	PHE	TRP	engineered mutation	UNP P0A251
D	81	PHE	TRP	engineered mutation	UNP P0A251
E	81	PHE	TRP	engineered mutation	UNP P0A251

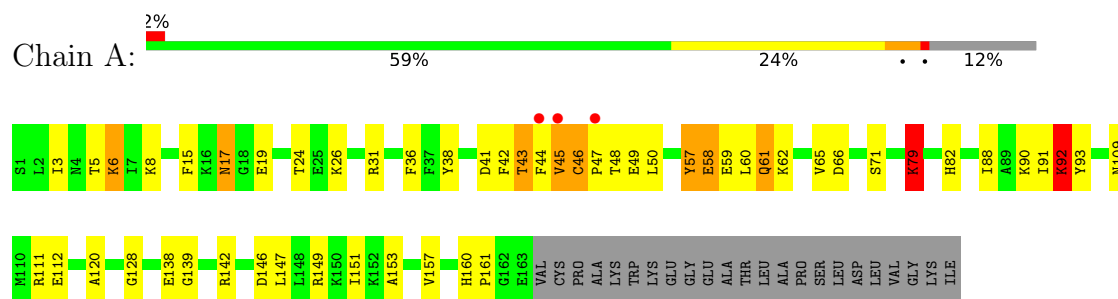
- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	K	0	0
			1	1		
2	B	2	Total	K	0	0
			2	2		
2	D	1	Total	K	0	0
			1	1		

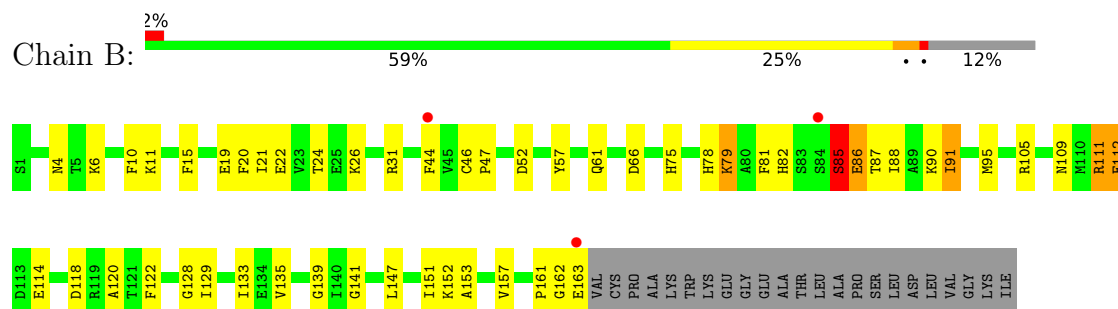
3 Residue-property plots [i](#)

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.

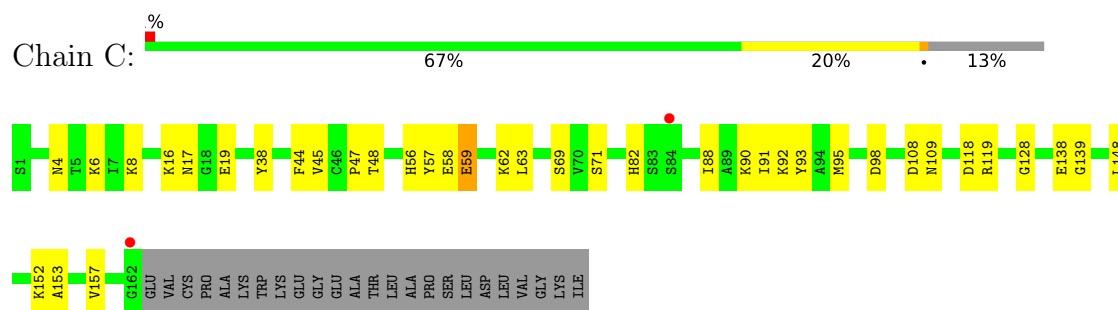
- Molecule 1: Alkyl hydroperoxide reductase subunit C



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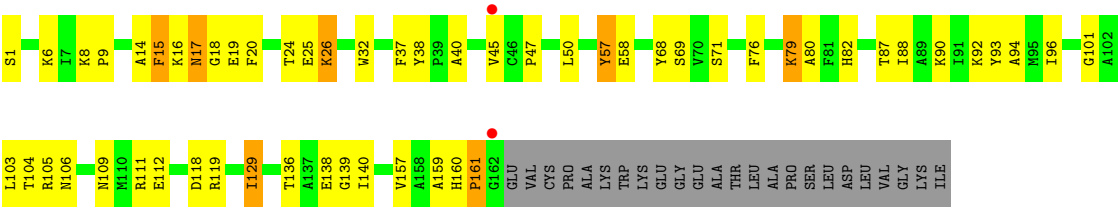


- Molecule 1: Alkyl hydroperoxide reductase subunit C

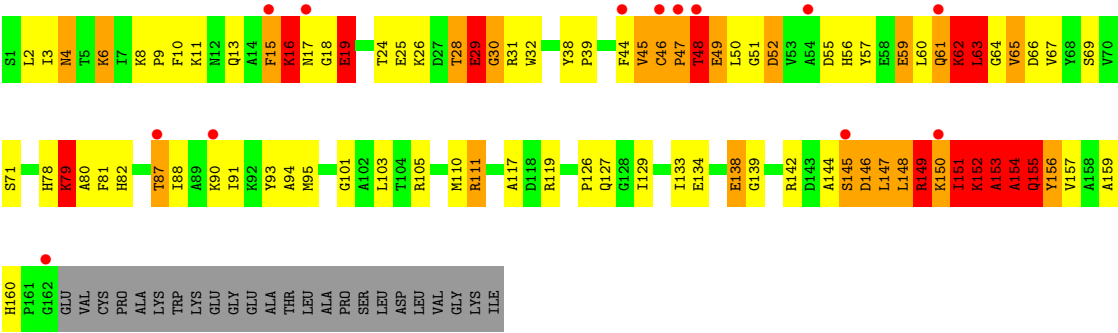


- Molecule 1: Alkyl hydroperoxide reductase subunit C





● Molecule 1: Alkyl hydroperoxide reductase subunit C



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	125.47Å 169.92Å 134.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.89 – 3.35 71.89 – 3.35	Depositor EDS
% Data completeness (in resolution range)	100.0 (71.89-3.35) 100.0 (71.89-3.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 3.33Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.235 , 0.279 0.230 , 0.276	Depositor DCC
R_{free} test set	1006 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	75.2	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 64.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6410	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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: K

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	0.99	0/1320	1.56	26/1787 (1.5%)
1	B	0.95	0/1314	1.41	15/1780 (0.8%)
1	C	0.87	1/1303 (0.1%)	1.43	13/1764 (0.7%)
1	D	2.10	7/1311 (0.5%)	1.38	15/1775 (0.8%)
1	E	2.28	17/1303 (1.3%)	2.20	71/1764 (4.0%)
All	All	1.56	25/6551 (0.4%)	1.62	140/8870 (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	A	0	2
1	E	0	7
All	All	0	9

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	79	LYS	CD-CE	54.13	3.14	1.52
1	D	79	LYS	CD-CE	51.97	3.08	1.52
1	D	15	PHE	CD1-CE1	22.53	2.06	1.38
1	E	15	PHE	CD1-CE1	20.30	1.99	1.38
1	D	15	PHE	CE1-CZ	19.49	1.97	1.38
1	D	15	PHE	CD2-CE2	18.84	1.95	1.38
1	D	15	PHE	CE2-CZ	18.55	1.94	1.38
1	E	15	PHE	CE1-CZ	18.17	1.93	1.38
1	E	15	PHE	CE2-CZ	17.46	1.91	1.38
1	E	15	PHE	CD2-CE2	17.28	1.90	1.38
1	E	15	PHE	CG-CD2	15.93	1.72	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	15	PHE	CG-CD1	15.33	1.71	1.38
1	E	15	PHE	CG-CD1	14.81	1.70	1.38
1	D	15	PHE	CG-CD2	13.74	1.67	1.38
1	E	65	VAL	CA-CB	10.89	1.67	1.54
1	E	153	ALA	N-CA	-9.48	1.33	1.46
1	E	150	LYS	N-CA	-9.04	1.35	1.46
1	E	152	LYS	CA-C	-8.60	1.41	1.52
1	E	48	THR	CA-CB	7.75	1.66	1.53
1	C	59	GLU	CA-C	7.02	1.61	1.52
1	E	16	LYS	CA-C	-6.49	1.45	1.52
1	E	13	GLN	CA-C	-5.43	1.45	1.52
1	E	48	THR	CA-C	5.24	1.59	1.52
1	E	65	VAL	CA-C	-5.03	1.46	1.52
1	E	133	ILE	C-O	-5.00	1.19	1.24

All (140) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	151	ILE	N-CA-C	-19.59	90.94	110.72
1	E	149	ARG	CA-C-N	-16.41	100.60	122.42
1	E	149	ARG	C-N-CA	-16.41	100.60	122.42
1	E	49	GLU	N-CA-C	-14.94	95.00	111.28
1	A	45	VAL	N-CA-C	-14.30	87.14	108.87
1	A	43	THR	N-CA-C	-14.24	88.29	110.36
1	E	79	LYS	CG-CD-CE	12.05	139.02	111.30
1	E	65	VAL	N-CA-CB	11.84	127.35	111.25
1	E	45	VAL	N-CA-C	-11.32	91.66	108.87
1	E	148	LEU	N-CA-C	-10.94	98.10	111.33
1	E	80	ALA	N-CA-C	-10.71	99.61	111.07
1	E	154	ALA	N-CA-C	-10.05	100.32	111.28
1	C	48	THR	N-CA-C	-9.91	100.48	111.28
1	E	46	CYS	CA-C-N	9.29	131.46	119.84
1	E	46	CYS	C-N-CA	9.29	131.46	119.84
1	E	29	GLU	CA-C-N	-9.21	105.64	121.85
1	E	29	GLU	C-N-CA	-9.21	105.64	121.85
1	E	65	VAL	CB-CA-C	-8.90	97.57	110.83
1	A	8	LYS	CA-C-N	8.83	128.88	119.78
1	A	8	LYS	C-N-CA	8.83	128.88	119.78
1	E	45	VAL	CA-C-N	-8.76	108.93	121.20
1	E	45	VAL	C-N-CA	-8.76	108.93	121.20
1	D	79	LYS	CG-CD-CE	8.63	131.14	111.30
1	E	46	CYS	N-CA-C	-8.45	96.92	109.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	85	SER	N-CA-C	8.43	122.99	109.24
1	E	61	GLN	N-CA-C	-8.29	102.18	111.71
1	C	8	LYS	CA-C-N	8.14	128.07	119.85
1	C	8	LYS	C-N-CA	8.14	128.07	119.85
1	A	46	CYS	CA-C-N	-8.03	109.80	119.84
1	A	46	CYS	C-N-CA	-8.03	109.80	119.84
1	E	17	ASN	N-CA-C	7.98	122.43	111.17
1	B	162	GLY	N-CA-C	7.87	131.84	113.18
1	C	17	ASN	CB-CA-C	-7.83	100.78	111.89
1	E	49	GLU	N-CA-CB	7.79	121.57	110.12
1	E	8	LYS	CA-C-N	7.79	127.80	119.78
1	E	8	LYS	C-N-CA	7.79	127.80	119.78
1	C	45	VAL	N-CA-C	-7.67	97.21	108.87
1	A	57	TYR	N-CA-C	7.67	119.28	111.07
1	D	45	VAL	N-CA-C	-7.43	99.10	109.21
1	A	59	GLU	N-CA-C	-7.35	103.27	111.28
1	E	62	LYS	CA-CB-CG	7.30	128.69	114.10
1	D	80	ALA	N-CA-C	-7.28	103.28	111.07
1	E	111	ARG	N-CA-C	-7.25	95.24	107.99
1	A	43	THR	CA-C-N	-7.23	110.83	122.62
1	A	43	THR	C-N-CA	-7.23	110.83	122.62
1	B	46	CYS	CA-C-N	7.21	128.85	119.84
1	B	46	CYS	C-N-CA	7.21	128.85	119.84
1	D	76	PHE	N-CA-C	-7.14	103.58	111.36
1	E	149	ARG	N-CA-CB	7.09	119.93	110.29
1	E	153	ALA	N-CA-C	-7.03	95.83	110.80
1	E	19	GLU	CA-C-N	7.02	132.09	122.19
1	E	19	GLU	C-N-CA	7.02	132.09	122.19
1	E	79	LYS	CB-CG-CD	-6.89	95.44	111.30
1	C	44	PHE	CA-C-N	-6.89	111.47	122.50
1	C	44	PHE	C-N-CA	-6.89	111.47	122.50
1	E	62	LYS	CA-C-N	6.82	135.95	122.31
1	E	62	LYS	C-N-CA	6.82	135.95	122.31
1	C	58	GLU	N-CA-C	-6.81	102.71	111.02
1	E	30	GLY	N-CA-C	-6.79	105.85	115.43
1	E	156	TYR	CA-CB-CG	6.72	126.00	113.90
1	A	57	TYR	CA-C-N	-6.71	109.39	121.14
1	A	57	TYR	C-N-CA	-6.71	109.39	121.14
1	E	147	LEU	N-CA-C	-6.71	103.89	111.07
1	A	41	ASP	N-CA-C	-6.65	101.57	110.55
1	D	79	LYS	CA-CB-CG	-6.59	100.92	114.10
1	D	17	ASN	N-CA-C	6.57	119.24	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	19	GLU	CA-CB-CG	-6.46	101.19	114.10
1	A	112	GLU	N-CA-C	6.42	119.09	111.71
1	C	17	ASN	N-CA-C	6.38	120.19	111.39
1	A	8	LYS	N-CA-C	-6.37	101.83	110.36
1	E	28	THR	CA-C-N	-6.33	113.55	124.82
1	E	28	THR	C-N-CA	-6.33	113.55	124.82
1	A	92	LYS	CB-CG-CD	6.32	125.83	111.30
1	E	47	PRO	CA-C-O	-6.31	109.12	120.60
1	E	91	ILE	N-CA-C	6.30	117.16	108.84
1	A	19	GLU	N-CA-C	6.29	119.79	109.85
1	B	86	GLU	CA-C-N	6.29	129.22	120.29
1	B	86	GLU	C-N-CA	6.29	129.22	120.29
1	A	61	GLN	CA-C-N	-6.18	112.78	122.49
1	A	61	GLN	C-N-CA	-6.18	112.78	122.49
1	E	138	GLU	CA-CB-CG	6.15	126.41	114.10
1	E	152	LYS	O-C-N	6.10	130.70	122.59
1	A	92	LYS	CA-CB-CG	-6.04	102.02	114.10
1	E	79	LYS	CD-CE-NZ	6.01	131.14	111.90
1	B	91	ILE	N-CA-C	5.90	116.63	108.84
1	E	18	GLY	N-CA-C	5.90	122.88	115.21
1	E	6	LYS	N-CA-C	-5.85	99.76	109.46
1	D	58	GLU	N-CA-CB	5.84	118.71	110.12
1	E	44	PHE	CA-C-N	-5.84	113.16	122.50
1	E	44	PHE	C-N-CA	-5.84	113.16	122.50
1	D	32	TRP	N-CA-C	-5.78	100.29	109.76
1	E	26	LYS	CA-CB-CG	-5.76	102.58	114.10
1	E	4	ASN	N-CA-C	5.71	119.55	112.24
1	E	65	VAL	CA-CB-CG2	5.71	120.10	110.40
1	E	52	ASP	N-CA-C	-5.69	105.08	111.28
1	A	61	GLN	N-CA-C	5.68	118.25	111.71
1	C	98	ASP	CA-C-N	5.66	126.04	119.47
1	C	98	ASP	C-N-CA	5.66	126.04	119.47
1	E	139	GLY	N-CA-C	-5.62	107.28	114.37
1	E	87	THR	N-CA-C	-5.61	105.24	111.36
1	B	52	ASP	N-CA-C	-5.61	105.17	111.28
1	E	152	LYS	CA-C-O	5.60	128.52	120.51
1	D	40	ALA	N-CA-C	5.59	117.59	108.76
1	D	19	GLU	N-CA-C	5.58	118.34	109.81
1	A	45	VAL	N-CA-CB	5.56	118.23	110.56
1	E	150	LYS	CA-C-N	5.55	128.31	120.42
1	E	150	LYS	C-N-CA	5.55	128.31	120.42
1	B	112	GLU	N-CA-C	5.52	118.06	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	79	LYS	CB-CA-C	-5.52	102.22	110.88
1	D	129	ILE	CB-CA-C	-5.51	103.99	111.15
1	E	62	LYS	O-C-N	5.51	129.91	122.59
1	B	46	CYS	CB-CA-C	5.48	118.22	109.62
1	E	152	LYS	CA-CB-CG	5.47	125.04	114.10
1	E	6	LYS	CA-CB-CG	5.43	124.97	114.10
1	D	57	TYR	CA-C-N	-5.35	113.12	120.28
1	D	57	TYR	C-N-CA	-5.35	113.12	120.28
1	B	57	TYR	N-CA-C	5.34	116.79	111.07
1	E	29	GLU	N-CA-CB	-5.33	103.61	113.89
1	E	29	GLU	N-CA-C	5.33	124.17	113.31
1	A	79	LYS	N-CA-CB	5.29	117.89	110.12
1	E	44	PHE	CB-CA-C	-5.28	101.63	110.56
1	B	44	PHE	CA-C-N	-5.28	114.05	122.50
1	B	44	PHE	C-N-CA	-5.28	114.05	122.50
1	A	58	GLU	CA-CB-CG	-5.28	103.55	114.10
1	C	90	LYS	CB-CG-CD	-5.25	99.21	111.30
1	D	111	ARG	N-CA-C	-5.23	98.15	107.98
1	E	150	LYS	CA-C-O	5.16	126.10	120.58
1	E	145	SER	CA-C-N	-5.12	113.45	123.13
1	E	145	SER	C-N-CA	-5.12	113.45	123.13
1	C	152	LYS	N-CA-C	-5.10	105.80	111.36
1	E	46	CYS	CA-C-O	-5.10	115.53	119.46
1	A	17	ASN	N-CA-CB	-5.09	104.16	111.70
1	E	44	PHE	N-CA-C	5.09	119.19	112.92
1	D	79	LYS	CD-CE-NZ	5.05	128.06	111.90
1	E	79	LYS	N-CA-C	5.05	116.78	111.28
1	A	46	CYS	N-CA-C	5.04	116.74	109.48
1	E	63	LEU	CB-CG-CD2	5.04	125.80	110.70
1	E	127	GLN	N-CA-C	-5.03	106.73	112.92
1	E	45	VAL	CB-CA-C	5.03	118.10	111.21
1	B	111	ARG	N-CA-C	-5.03	99.14	107.99

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	45	VAL	Peptide
1	A	62	LYS	Peptide
1	E	146	ASP	Peptide
1	E	149	ARG	Peptide
1	E	151	ILE	Peptide

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Mol	Chain	Res	Type	Group
1	E	154	ALA	Peptide
1	E	19	GLU	Peptide
1	E	28	THR	Peptide
1	E	29	GLU	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	A	1291	0	1243	41	0
1	B	1285	0	1230	39	0
1	C	1274	0	1232	25	0
1	D	1282	0	1237	75	0
1	E	1274	0	1232	151	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	D	1	0	0	0	0
All	All	6410	0	6174	324	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 26.

All (324) 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:E:15:PHE:CE1	1:E:15:PHE:CZ	1.93	1.57
1:D:15:PHE:CE2	1:D:15:PHE:CZ	1.94	1.56
1:E:15:PHE:CZ	1:E:15:PHE:CE2	1.91	1.55
1:E:15:PHE:CE2	1:E:15:PHE:CD2	1.90	1.54
1:D:15:PHE:CZ	1:D:15:PHE:CE1	1.97	1.53
1:D:15:PHE:CE2	1:D:15:PHE:CD2	1.95	1.50
1:E:15:PHE:CE1	1:E:15:PHE:CD1	1.99	1.50
1:D:15:PHE:CE1	1:D:15:PHE:CD1	2.06	1.41
1:E:29:GLU:HG2	1:E:31:ARG:HG3	1.24	1.09
1:D:15:PHE:CD1	1:D:79:LYS:CE	2.35	1.08
1:E:63:LEU:HD11	1:E:65:VAL:HG23	1.30	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:15:PHE:CE1	1:D:79:LYS:CD	2.36	1.08
1:E:15:PHE:CZ	1:E:79:LYS:CE	2.39	1.06
1:E:146:ASP:HB2	1:E:149:ARG:HB2	1.35	1.06
1:E:15:PHE:CD1	1:E:79:LYS:CD	2.40	1.05
1:D:15:PHE:CG	1:D:79:LYS:CE	2.39	1.05
1:E:15:PHE:CD2	1:E:79:LYS:CD	2.40	1.05
1:E:15:PHE:CE2	1:E:79:LYS:CE	2.39	1.04
1:D:15:PHE:CD2	1:D:79:LYS:CE	2.40	1.04
1:D:15:PHE:CZ	1:D:79:LYS:CD	2.42	1.02
1:D:15:PHE:CE1	1:D:79:LYS:CE	2.42	1.02
1:D:15:PHE:CE2	1:D:79:LYS:CE	2.42	1.02
1:E:15:PHE:CD1	1:E:79:LYS:HD3	1.94	1.02
1:E:15:PHE:CE1	1:E:79:LYS:CE	2.43	1.02
1:E:146:ASP:HA	1:E:149:ARG:H	1.19	1.01
1:D:15:PHE:CG	1:D:79:LYS:HE3	1.96	1.01
1:E:15:PHE:CE2	1:E:79:LYS:CD	2.44	1.01
1:E:15:PHE:CG	1:E:79:LYS:CE	2.44	1.01
1:E:15:PHE:CG	1:E:79:LYS:CD	2.44	1.00
1:E:15:PHE:CD2	1:E:79:LYS:HD2	1.96	0.99
1:E:15:PHE:CD2	1:E:79:LYS:CE	2.44	0.99
1:E:15:PHE:CE1	1:E:79:LYS:CD	2.45	0.99
1:D:15:PHE:CD1	1:D:79:LYS:CD	2.45	0.99
1:C:59:GLU:O	1:C:62:LYS:HG2	1.63	0.98
1:D:15:PHE:CZ	1:D:79:LYS:CE	2.46	0.98
1:D:15:PHE:CE2	1:D:79:LYS:CD	2.47	0.97
1:E:15:PHE:CZ	1:E:79:LYS:HE3	1.96	0.97
1:E:15:PHE:CZ	1:E:79:LYS:CD	2.49	0.96
1:B:15:PHE:HD1	1:B:79:LYS:HZ2	0.98	0.96
1:E:15:PHE:CD1	1:E:79:LYS:CE	2.48	0.96
1:D:15:PHE:CE2	1:D:79:LYS:HE2	2.02	0.95
1:C:119:ARG:HE	1:C:138:GLU:HG2	1.31	0.93
1:D:15:PHE:CD2	1:D:79:LYS:CD	2.52	0.92
1:E:29:GLU:HG2	1:E:31:ARG:CG	1.98	0.92
1:D:15:PHE:CG	1:D:79:LYS:CD	2.55	0.89
1:E:150:LYS:C	1:E:152:LYS:HB2	1.96	0.89
1:A:146:ASP:OD1	1:A:149:ARG:NH1	2.06	0.88
1:E:152:LYS:HB3	1:E:154:ALA:N	1.88	0.88
1:E:146:ASP:HB2	1:E:149:ARG:CB	2.03	0.87
1:E:152:LYS:HB3	1:E:154:ALA:H	1.37	0.87
1:D:15:PHE:CZ	1:D:79:LYS:HD2	2.05	0.87
1:E:63:LEU:HD11	1:E:65:VAL:CG2	2.04	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:59:GLU:HG3	1:C:62:LYS:HE3	1.58	0.85
1:E:29:GLU:CG	1:E:31:ARG:HG3	2.06	0.84
1:E:150:LYS:O	1:E:152:LYS:HB2	1.78	0.84
1:E:153:ALA:HB1	1:E:156:TYR:CB	2.07	0.84
1:E:15:PHE:CD1	1:E:79:LYS:HE2	2.14	0.83
1:E:59:GLU:O	1:E:62:LYS:HB2	1.78	0.82
1:E:16:LYS:HA	1:E:79:LYS:HZ3	1.45	0.82
1:D:15:PHE:CE2	1:D:79:LYS:HD3	2.15	0.79
1:A:46:CYS:HB3	1:A:48:THR:HG23	1.64	0.79
1:E:63:LEU:CD1	1:E:65:VAL:HG23	2.10	0.79
1:D:15:PHE:CE1	1:D:79:LYS:HD2	2.16	0.78
1:E:62:LYS:NZ	1:E:63:LEU:HB3	1.97	0.78
1:D:15:PHE:CD2	1:D:79:LYS:HE3	2.16	0.78
1:E:153:ALA:HB1	1:E:156:TYR:HB3	1.66	0.77
1:B:78:HIS:HD2	1:B:95:MET:HE2	1.47	0.77
1:B:15:PHE:HD1	1:B:79:LYS:NZ	1.82	0.76
1:D:17:ASN:N	1:D:79:LYS:HZ1	1.83	0.76
1:A:157:VAL:HG11	1:B:139:GLY:HA3	1.67	0.76
1:E:62:LYS:HB3	1:E:63:LEU:HD23	1.68	0.76
1:E:146:ASP:CB	1:E:149:ARG:HB2	2.13	0.75
1:E:57:TYR:CE1	1:E:93:TYR:HB3	2.22	0.75
1:C:56:HIS:HD2	1:C:148:LEU:HD12	1.51	0.74
1:B:15:PHE:CD1	1:B:79:LYS:HD3	2.23	0.74
1:E:62:LYS:HZ2	1:E:63:LEU:HB3	1.50	0.74
1:E:31:ARG:NH1	1:E:66:ASP:OD1	2.20	0.73
1:D:101:GLY:O	1:D:105:ARG:HG3	1.87	0.73
1:E:146:ASP:HA	1:E:149:ARG:N	2.02	0.73
1:A:139:GLY:HA3	1:B:157:VAL:HG11	1.71	0.73
1:E:31:ARG:NH2	1:E:64:GLY:HA2	2.03	0.72
1:D:15:PHE:CD2	1:D:79:LYS:HD3	2.24	0.72
1:E:148:LEU:C	1:E:150:LYS:HZ2	1.97	0.72
1:E:119:ARG:HE	1:E:138:GLU:HB3	1.56	0.70
1:E:101:GLY:O	1:E:105:ARG:HG2	1.91	0.70
1:B:152:LYS:HG2	1:B:153:ALA:N	2.05	0.70
1:D:15:PHE:CE1	1:D:79:LYS:NZ	2.59	0.69
1:D:82:HIS:HA	1:D:88:ILE:HG22	1.75	0.68
1:B:88:ILE:HG23	1:B:91:ILE:HD12	1.74	0.68
1:D:18:GLY:N	1:D:79:LYS:NZ	2.41	0.68
1:E:16:LYS:CA	1:E:79:LYS:HZ3	2.07	0.67
1:E:63:LEU:HD13	1:E:155:GLN:HE22	1.58	0.67
1:E:15:PHE:CE1	1:E:79:LYS:HE2	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:47:PRO:C	1:E:49:GLU:H	2.03	0.66
1:D:15:PHE:CD1	1:D:79:LYS:CG	2.79	0.66
1:E:152:LYS:HZ2	1:E:154:ALA:HA	1.61	0.65
1:C:62:LYS:HG3	1:C:63:LEU:N	2.10	0.65
1:A:15:PHE:CE1	1:A:79:LYS:HB2	2.32	0.65
1:B:20:PHE:CE1	1:B:79:LYS:HE2	2.31	0.65
1:A:6:LYS:HA	1:A:128:GLY:O	1.96	0.65
1:B:109:ASN:OD1	1:B:118:ASP:HB2	1.97	0.64
1:E:56:HIS:O	1:E:59:GLU:HG2	1.98	0.64
1:A:17:ASN:ND2	1:A:92:LYS:HG2	2.13	0.64
1:E:150:LYS:HB3	1:E:152:LYS:HA	1.79	0.64
1:B:87:THR:O	1:B:90:LYS:HB2	1.98	0.63
1:B:153:ALA:O	1:B:157:VAL:HG22	1.99	0.63
1:E:62:LYS:HZ2	1:E:63:LEU:CB	2.11	0.63
1:C:108:ASP:O	1:D:1:SER:N	2.30	0.63
1:E:50:LEU:O	1:E:93:TYR:OH	2.16	0.63
1:E:62:LYS:HD3	1:E:63:LEU:CD2	2.29	0.63
1:E:47:PRO:O	1:E:49:GLU:N	2.32	0.63
1:C:6:LYS:HA	1:C:128:GLY:O	1.98	0.62
1:D:119:ARG:HE	1:D:138:GLU:HA	1.65	0.62
1:E:32:TRP:HB3	1:E:65:VAL:HG22	1.82	0.62
1:D:15:PHE:CZ	1:D:79:LYS:HE2	2.34	0.61
1:E:152:LYS:NZ	1:E:154:ALA:HA	2.15	0.61
1:A:17:ASN:ND2	1:A:92:LYS:CG	2.64	0.61
1:D:109:ASN:OD1	1:D:118:ASP:HB2	2.01	0.61
1:E:78:HIS:HD2	1:E:95:MET:HB3	1.66	0.61
1:E:87:THR:O	1:E:90:LYS:HB2	2.01	0.60
1:D:18:GLY:N	1:D:79:LYS:HZ3	1.99	0.60
1:D:15:PHE:HE2	1:D:82:HIS:CG	2.18	0.60
1:E:153:ALA:HB1	1:E:156:TYR:HB2	1.81	0.60
1:E:15:PHE:CD2	1:E:79:LYS:NZ	2.70	0.60
1:E:16:LYS:NZ	1:E:94:ALA:HB2	2.17	0.60
1:E:78:HIS:CD2	1:E:95:MET:HB3	2.36	0.60
1:E:15:PHE:CE2	1:E:79:LYS:HE3	2.36	0.59
1:B:78:HIS:CD2	1:B:95:MET:HE2	2.33	0.59
1:E:156:TYR:CD1	1:E:160:HIS:HB2	2.38	0.59
1:C:59:GLU:HG3	1:C:62:LYS:CE	2.29	0.59
1:E:152:LYS:HG2	1:E:153:ALA:HB3	1.85	0.58
1:E:150:LYS:HB3	1:E:152:LYS:CA	2.33	0.58
1:A:17:ASN:HD21	1:A:92:LYS:HG2	1.68	0.58
1:B:15:PHE:HE2	1:B:82:HIS:CD2	2.21	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:156:TYR:O	1:E:160:HIS:N	2.36	0.57
1:A:111:ARG:HH22	1:A:138:GLU:HG3	1.68	0.57
1:A:58:GLU:HA	1:A:61:GLN:HB2	1.85	0.57
1:E:52:ASP:OD1	1:E:144:ALA:HB3	2.04	0.57
1:B:31:ARG:NH1	1:B:66:ASP:OD1	2.38	0.56
1:E:149:ARG:HA	1:E:150:LYS:HG2	1.86	0.56
1:E:149:ARG:HG2	1:E:150:LYS:N	2.20	0.56
1:E:48:THR:HA	1:E:51:GLY:H	1.71	0.56
1:A:149:ARG:NH2	1:B:141:GLY:O	2.39	0.56
1:A:31:ARG:NH1	1:A:66:ASP:OD1	2.40	0.55
1:E:15:PHE:HE2	1:E:82:HIS:CB	2.20	0.55
1:D:16:LYS:C	1:D:79:LYS:HZ1	2.15	0.55
1:E:63:LEU:HD21	1:E:65:VAL:HG23	1.89	0.55
1:D:105:ARG:HH21	1:D:112:GLU:CD	2.15	0.55
1:B:81:PHE:O	1:B:85:SER:OG	2.25	0.55
1:E:134:GLU:OE1	1:E:142:ARG:HG2	2.07	0.55
1:C:109:ASN:OD1	1:C:118:ASP:HB2	2.08	0.54
1:D:38:TYR:CD2	1:D:50:LEU:HD21	2.42	0.54
1:D:82:HIS:HA	1:D:88:ILE:CG2	2.37	0.54
1:B:6:LYS:HA	1:B:128:GLY:O	2.07	0.54
1:D:9:PRO:HA	1:D:25:GLU:HB3	1.89	0.54
1:E:10:PHE:O	1:E:24:THR:HA	2.06	0.54
1:E:16:LYS:HZ1	1:E:94:ALA:HB2	1.73	0.54
1:E:144:ALA:HA	1:E:147:LEU:HB3	1.90	0.54
1:C:62:LYS:HG3	1:C:63:LEU:HG	1.90	0.54
1:D:87:THR:O	1:D:90:LYS:HG3	2.08	0.54
1:C:82:HIS:HA	1:C:88:ILE:HG22	1.90	0.53
1:E:149:ARG:CG	1:E:150:LYS:N	2.71	0.53
1:B:31:ARG:HD3	1:B:66:ASP:OD2	2.08	0.53
1:D:16:LYS:C	1:D:79:LYS:NZ	2.67	0.53
1:E:32:TRP:O	1:E:65:VAL:HA	2.09	0.53
1:D:8:LYS:HE2	1:D:106:ASN:OD1	2.09	0.53
1:A:157:VAL:HG11	1:B:139:GLY:CA	2.37	0.53
1:B:147:LEU:O	1:B:151:ILE:HG13	2.09	0.53
1:E:61:GLN:O	1:E:64:GLY:N	2.30	0.53
1:A:58:GLU:HG2	1:A:61:GLN:HB2	1.90	0.52
1:E:47:PRO:C	1:E:49:GLU:N	2.67	0.52
1:B:15:PHE:CD1	1:B:79:LYS:CD	2.92	0.52
1:E:105:ARG:HH21	1:E:110:MET:HG2	1.75	0.52
1:C:88:ILE:HG23	1:C:91:ILE:HD12	1.91	0.52
1:E:59:GLU:CD	1:E:59:GLU:H	2.18	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:146:ASP:CA	1:E:149:ARG:H	2.08	0.51
1:A:109:ASN:ND2	1:A:120:ALA:HB3	2.26	0.51
1:E:32:TRP:CB	1:E:65:VAL:HG22	2.39	0.51
1:D:105:ARG:NH2	1:D:112:GLU:OE1	2.44	0.51
1:E:57:TYR:CZ	1:E:61:GLN:NE2	2.78	0.51
1:E:61:GLN:O	1:E:62:LYS:C	2.53	0.51
1:B:47:PRO:HB2	1:B:88:ILE:HD11	1.92	0.51
1:B:105:ARG:NH2	1:B:112:GLU:OE1	2.43	0.51
1:B:78:HIS:CD2	1:B:95:MET:HB3	2.46	0.51
1:C:157:VAL:HG11	1:D:139:GLY:HA3	1.93	0.51
1:B:11:LYS:HZ2	1:B:22:GLU:CD	2.18	0.51
1:C:59:GLU:HA	1:C:62:LYS:HE3	1.93	0.51
1:D:18:GLY:N	1:D:79:LYS:HZ1	2.08	0.51
1:E:150:LYS:HB3	1:E:152:LYS:CB	2.40	0.50
1:A:38:TYR:CZ	1:A:71:SER:HB3	2.46	0.50
1:A:147:LEU:O	1:A:151:ILE:HG13	2.11	0.50
1:C:47:PRO:CB	1:C:88:ILE:HD11	2.42	0.50
1:A:36:PHE:HZ	1:A:49:GLU:HB3	1.77	0.50
1:A:65:VAL:HG21	1:A:151:ILE:HG21	1.93	0.49
1:B:15:PHE:CE2	1:B:82:HIS:CG	2.99	0.49
1:C:38:TYR:CZ	1:C:71:SER:HB3	2.47	0.49
1:B:163:GLU:N	1:B:163:GLU:OE1	2.45	0.49
1:E:146:ASP:O	1:E:149:ARG:HB3	2.12	0.49
1:B:111:ARG:HD3	1:B:114:GLU:OE1	2.13	0.49
1:B:10:PHE:O	1:B:24:THR:HA	2.12	0.49
1:E:15:PHE:CE1	1:E:79:LYS:HD3	2.43	0.49
1:E:153:ALA:CB	1:E:156:TYR:HB3	2.40	0.49
1:D:17:ASN:CA	1:D:79:LYS:HZ1	2.26	0.49
1:E:39:PRO:HG3	1:E:111:ARG:NH2	2.28	0.49
1:A:24:THR:OG1	1:A:26:LYS:HG2	2.13	0.49
1:E:119:ARG:HE	1:E:138:GLU:CB	2.24	0.49
1:D:16:LYS:HB2	1:D:94:ALA:HA	1.95	0.48
1:E:63:LEU:CD2	1:E:65:VAL:HG23	2.44	0.48
1:E:31:ARG:HD3	1:E:66:ASP:OD2	2.13	0.48
1:E:55:ASP:OD1	1:E:90:LYS:HD3	2.14	0.48
1:E:149:ARG:HA	1:E:150:LYS:CG	2.44	0.48
1:E:156:TYR:O	1:E:159:ALA:N	2.45	0.48
1:E:6:LYS:HA	1:E:129:ILE:HD13	1.96	0.47
1:B:11:LYS:NZ	1:B:22:GLU:OE1	2.47	0.47
1:E:38:TYR:CZ	1:E:71:SER:HB3	2.50	0.47
1:E:149:ARG:HA	1:E:150:LYS:NZ	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ASN:ND2	1:A:92:LYS:HG3	2.29	0.47
1:D:79:LYS:HD3	1:D:79:LYS:HA	1.39	0.47
1:E:15:PHE:CZ	1:E:79:LYS:CG	2.96	0.47
1:E:30:GLY:HA2	1:E:126:PRO:HB3	1.97	0.47
1:E:153:ALA:O	1:E:156:TYR:HB3	2.14	0.47
1:D:26:LYS:H	1:D:26:LYS:HG2	1.61	0.47
1:A:5:THR:HG22	1:A:6:LYS:N	2.30	0.47
1:A:50:LEU:HB3	1:A:91:ILE:HD11	1.96	0.47
1:A:49:GLU:OE2	1:A:142:ARG:N	2.44	0.46
1:E:150:LYS:CB	1:E:152:LYS:HA	2.44	0.46
1:A:82:HIS:HA	1:A:88:ILE:HG22	1.96	0.46
1:B:122:PHE:CD2	1:B:133:ILE:HG12	2.50	0.46
1:D:15:PHE:CD1	1:D:79:LYS:NZ	2.83	0.46
1:D:17:ASN:HD21	1:D:92:LYS:HG3	1.80	0.46
1:E:9:PRO:HA	1:E:25:GLU:HB3	1.98	0.46
1:A:31:ARG:HD3	1:A:66:ASP:OD2	2.15	0.46
1:C:57:TYR:HD2	1:C:92:LYS:HD3	1.79	0.46
1:E:156:TYR:O	1:E:157:VAL:C	2.59	0.46
1:E:62:LYS:HD3	1:E:63:LEU:HD22	1.98	0.46
1:E:145:SER:O	1:E:148:LEU:HB2	2.15	0.46
1:C:59:GLU:CG	1:C:62:LYS:HE3	2.37	0.46
1:E:57:TYR:CE2	1:E:61:GLN:NE2	2.84	0.46
1:E:62:LYS:HD3	1:E:63:LEU:HD23	1.98	0.46
1:E:156:TYR:CD1	1:E:160:HIS:CG	3.04	0.46
1:D:136:THR:HB	1:D:140:ILE:HB	1.97	0.45
1:B:15:PHE:HE2	1:B:82:HIS:CG	2.34	0.45
1:E:150:LYS:CD	1:E:152:LYS:HA	2.46	0.45
1:D:15:PHE:C	1:D:79:LYS:HE3	2.41	0.45
1:E:39:PRO:HA	1:E:117:ALA:O	2.17	0.45
1:E:156:TYR:CE1	1:E:160:HIS:HB2	2.51	0.45
1:E:69:SER:OG	1:E:95:MET:HA	2.17	0.45
1:E:153:ALA:C	1:E:156:TYR:H	2.24	0.45
1:A:3:ILE:HD12	1:A:3:ILE:HA	1.83	0.45
1:B:6:LYS:HA	1:B:129:ILE:HD13	1.99	0.45
1:E:31:ARG:HH21	1:E:64:GLY:HA2	1.78	0.45
1:C:69:SER:OG	1:C:95:MET:HA	2.17	0.45
1:C:91:ILE:HG23	1:C:93:TYR:CZ	2.52	0.44
1:D:159:ALA:C	1:D:161:PRO:HD3	2.41	0.44
1:E:62:LYS:CE	1:E:63:LEU:HB3	2.46	0.44
1:C:16:LYS:O	1:C:19:GLU:HG2	2.17	0.44
1:D:57:TYR:CZ	1:D:93:TYR:HB3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:62:LYS:CG	1:C:63:LEU:N	2.79	0.44
1:E:45:VAL:O	1:E:47:PRO:HD3	2.17	0.44
1:E:57:TYR:CZ	1:E:93:TYR:HB3	2.50	0.44
1:E:62:LYS:CB	1:E:63:LEU:HD23	2.42	0.44
1:E:153:ALA:O	1:E:156:TYR:N	2.49	0.44
1:D:17:ASN:HD21	1:D:92:LYS:CG	2.31	0.44
1:D:17:ASN:C	1:D:79:LYS:HZ1	2.25	0.44
1:E:49:GLU:OE1	1:E:119:ARG:NH2	2.50	0.44
1:E:149:ARG:N	1:E:150:LYS:HZ2	2.15	0.44
1:E:150:LYS:C	1:E:152:LYS:CB	2.80	0.44
1:A:17:ASN:HD21	1:A:92:LYS:CG	2.29	0.43
1:A:58:GLU:HG2	1:A:61:GLN:HG3	1.99	0.43
1:B:20:PHE:CE1	1:B:75:HIS:HD2	2.37	0.43
1:D:6:LYS:HE2	1:D:6:LYS:HB2	1.60	0.43
1:E:88:ILE:HD13	1:E:88:ILE:HA	1.83	0.43
1:D:15:PHE:CG	1:D:79:LYS:CG	3.02	0.43
1:E:153:ALA:HB1	1:E:156:TYR:CD2	2.53	0.43
1:E:153:ALA:CA	1:E:156:TYR:H	2.31	0.43
1:A:15:PHE:CD1	1:A:79:LYS:HB2	2.52	0.43
1:B:21:ILE:HG22	1:B:22:GLU:N	2.34	0.43
1:E:67:VAL:O	1:E:93:TYR:HB2	2.17	0.43
1:A:153:ALA:O	1:A:157:VAL:HG22	2.17	0.43
1:D:20:PHE:CE1	1:D:79:LYS:HG3	2.53	0.43
1:D:47:PRO:HB2	1:D:88:ILE:HD11	2.01	0.43
1:A:57:TYR:HA	1:A:60:LEU:HD12	2.00	0.43
1:D:38:TYR:HE2	1:D:69:SER:HB2	1.83	0.43
1:E:78:HIS:O	1:E:81:PHE:HB3	2.18	0.43
1:E:151:ILE:O	1:E:154:ALA:N	2.46	0.43
1:E:153:ALA:C	1:E:156:TYR:N	2.77	0.43
1:D:14:ALA:HB2	1:D:96:ILE:HG12	2.00	0.43
1:E:62:LYS:HB3	1:E:63:LEU:H	1.08	0.43
1:E:63:LEU:HB2	1:E:155:GLN:NE2	2.34	0.43
1:E:78:HIS:HD2	1:E:95:MET:HE2	1.84	0.43
1:A:44:PHE:CE1	1:A:47:PRO:HD3	2.55	0.42
1:D:16:LYS:N	1:D:79:LYS:HE3	2.35	0.42
1:A:42:PHE:C	1:A:43:THR:O	2.54	0.42
1:B:120:ALA:HB2	1:B:135:VAL:HG13	2.02	0.42
1:D:15:PHE:CD1	1:D:79:LYS:HG3	2.52	0.42
1:E:153:ALA:HA	1:E:156:TYR:H	1.83	0.42
1:C:56:HIS:CD2	1:C:148:LEU:HD12	2.43	0.42
1:D:129:ILE:HG23	1:D:129:ILE:HD12	1.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:146:ASP:CB	1:E:149:ARG:CB	2.86	0.42
1:C:139:GLY:HA3	1:D:157:VAL:HG11	2.02	0.42
1:A:91:ILE:HG23	1:A:93:TYR:CZ	2.54	0.42
1:B:61:GLN:HE21	1:B:61:GLN:HA	1.85	0.42
1:D:159:ALA:C	1:D:160:HIS:HD2	2.28	0.42
1:D:9:PRO:HB3	1:D:25:GLU:CD	2.45	0.42
1:B:20:PHE:CE1	1:B:79:LYS:CE	3.02	0.42
1:D:24:THR:OG1	1:D:26:LYS:HG2	2.20	0.42
1:E:147:LEU:O	1:E:149:ARG:O	2.38	0.42
1:D:18:GLY:H	1:D:79:LYS:NZ	2.16	0.41
1:E:30:GLY:C	1:E:126:PRO:HB3	2.45	0.41
1:A:24:THR:OG1	1:A:26:LYS:CG	2.68	0.41
1:E:2:LEU:O	1:E:3:ILE:C	2.64	0.41
1:E:57:TYR:O	1:E:60:LEU:N	2.54	0.41
1:A:44:PHE:CD2	1:A:44:PHE:O	2.74	0.41
1:C:153:ALA:O	1:C:157:VAL:HG22	2.21	0.41
1:A:160:HIS:N	1:A:161:PRO:HD3	2.36	0.40
1:D:37:PHE:CE2	1:D:104:THR:HA	2.56	0.40
1:E:15:PHE:CE1	1:E:79:LYS:HE3	2.48	0.40
1:E:150:LYS:C	1:E:152:LYS:N	2.74	0.40
1:A:46:CYS:C	1:A:48:THR:H	2.29	0.40
1:A:58:GLU:HG2	1:A:61:GLN:CB	2.52	0.40
1:D:68:TYR:CD1	1:D:96:ILE:HD11	2.57	0.40
1:D:68:TYR:HB3	1:D:96:ILE:HD12	2.02	0.40
1:E:62:LYS:HB3	1:E:63:LEU:CD2	2.43	0.40
1:E:105:ARG:NH2	1:E:110:MET:HE3	2.36	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	162/186 (87%)	156 (96%)	6 (4%)	0	100	100
1	B	162/186 (87%)	156 (96%)	4 (2%)	2 (1%)	10	33
1	C	160/186 (86%)	155 (97%)	5 (3%)	0	100	100
1	D	161/186 (87%)	156 (97%)	4 (2%)	1 (1%)	21	49
1	E	160/186 (86%)	147 (92%)	8 (5%)	5 (3%)	3	17
All	All	805/930 (87%)	770 (96%)	27 (3%)	8 (1%)	12	38

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	86	GLU
1	E	62	LYS
1	E	152	LYS
1	E	155	GLN
1	E	48	THR
1	E	153	ALA
1	B	161	PRO
1	D	161	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	137/154 (89%)	133 (97%)	4 (3%)	37	60
1	B	136/154 (88%)	132 (97%)	4 (3%)	37	60
1	C	135/154 (88%)	134 (99%)	1 (1%)	76	77
1	D	136/154 (88%)	133 (98%)	3 (2%)	45	63
1	E	135/154 (88%)	123 (91%)	12 (9%)	9	31
All	All	679/770 (88%)	655 (96%)	24 (4%)	31	57

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	79	LYS
1	A	90	LYS
1	A	92	LYS
1	B	4	ASN
1	B	19	GLU
1	B	26	LYS
1	B	85	SER
1	C	4	ASN
1	D	26	LYS
1	D	71	SER
1	D	103	LEU
1	E	4	ASN
1	E	11	LYS
1	E	16	LYS
1	E	19	GLU
1	E	46	CYS
1	E	48	THR
1	E	59	GLU
1	E	63	LEU
1	E	79	LYS
1	E	103	LEU
1	E	152	LYS
1	E	155	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	160	HIS
1	B	17	ASN
1	C	61	GLN
1	D	61	GLN
1	E	17	ASN
1	E	82	HIS
1	E	155	GLN

5.3.3 RNA ⓘ

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 4 ligands modelled in this entry, 4 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	A	163/186 (87%)	-0.29	3 (1%)	67	51	23, 62, 105, 169	1 (0%)
1	B	163/186 (87%)	-0.16	3 (1%)	67	51	24, 69, 112, 166	1 (0%)
1	C	162/186 (87%)	-0.13	2 (1%)	76	63	45, 74, 118, 142	0
1	D	162/186 (87%)	0.01	2 (1%)	76	63	40, 84, 125, 169	1 (0%)
1	E	162/186 (87%)	0.48	13 (8%)	18	15	73, 114, 168, 213	0
All	All	812/930 (87%)	-0.02	23 (2%)	55	40	23, 80, 144, 213	3 (0%)

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	44	PHE	4.9
1	E	47	PRO	4.2
1	E	162	GLY	4.2
1	B	84	SER	3.7
1	E	44	PHE	3.7
1	A	47	PRO	3.1
1	A	44	PHE	3.0
1	D	162	GLY	2.9
1	D	45	VAL	2.8
1	C	84	SER	2.7
1	E	150	LYS	2.6
1	A	45	VAL	2.6
1	E	87	THR	2.5
1	C	162	GLY	2.4
1	E	90	LYS	2.4
1	E	48	THR	2.2
1	E	145	SER	2.2
1	E	61	GLN	2.2
1	E	46	CYS	2.2
1	E	17	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	163	GLU	2.1
1	E	15	PHE	2.1
1	E	54	ALA	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](#)

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(Å ²)	Q<0.9
2	K	D	201	1/1	0.87	0.12	118,118,118,118	0
2	K	B	202	1/1	0.89	0.07	71,71,71,71	0
2	K	B	201	1/1	0.94	0.07	83,83,83,83	0
2	K	A	201	1/1	0.97	0.20	48,48,48,48	1

6.5 Other polymers [i](#)

There are no such residues in this entry.