



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2024 – 11:47 PM EST

PDB ID : 2D4A
Title : Structure of the malate dehydrogenase from *Aeropyrum pernix*
Authors : Kawakami, R.; Sakuraba, H.; Tsuge, H.; Ohshima, T.
Deposited on : 2005-10-12
Resolution : 2.87 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

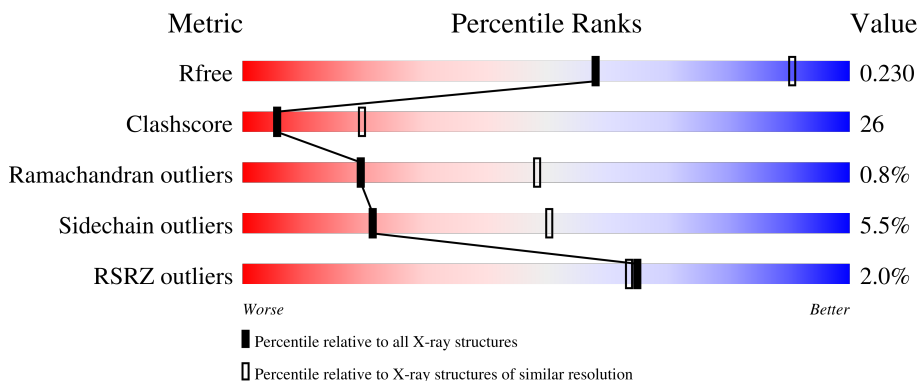
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.87 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}	130704	2691 (2.90-2.86)
Clashscore	141614	2947 (2.90-2.86)
Ramachandran outliers	138981	2868 (2.90-2.86)
Sidechain outliers	138945	2871 (2.90-2.86)
RSRZ outliers	127900	2629 (2.90-2.86)

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	308	 2% 56% 35% 5% ••
1	B	308	 2% 61% 30% 7% •
1	C	308	 2% 50% 39% 7% ••
1	D	308	 3% 59% 36% 5% •

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 9292 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 Malate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	301	2289	1455	383	438	13	0	0	0
1	D	308	2346	1489	397	446	14	0	0	0
1	C	299	2274	1446	384	431	13	0	0	0
1	A	298	2262	1440	379	430	13	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q9YEA1
B	1	MET	-	initiating methionine	UNP Q9YEA1
C	1	MET	-	initiating methionine	UNP Q9YEA1
D	1	MET	-	initiating methionine	UNP Q9YEA1

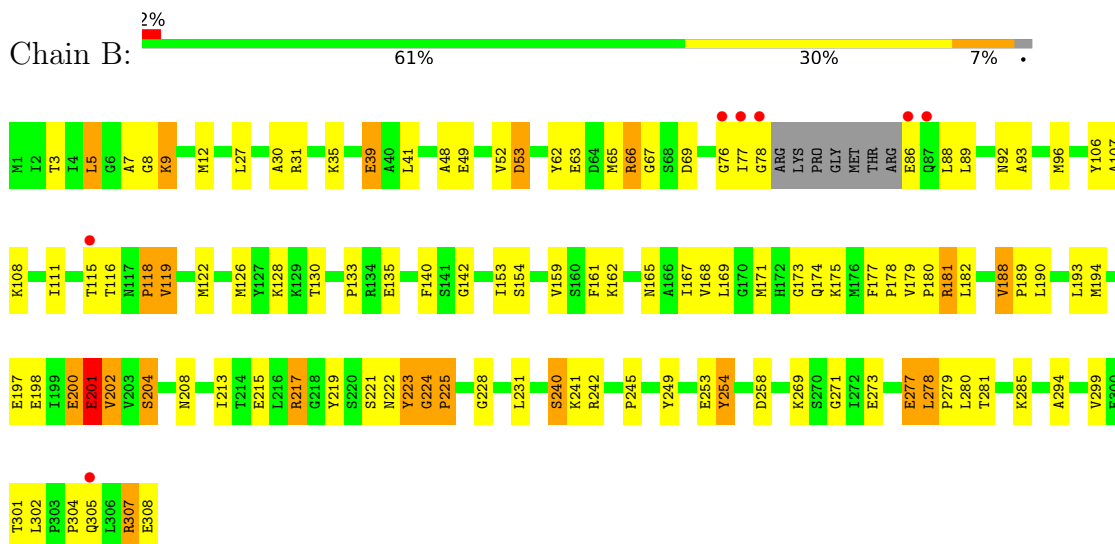
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	42	Total	O	0	0
			42	42		
2	D	34	Total	O	0	0
			34	34		
2	C	17	Total	O	0	0
			17	17		
2	A	28	Total	O	0	0
			28	28		

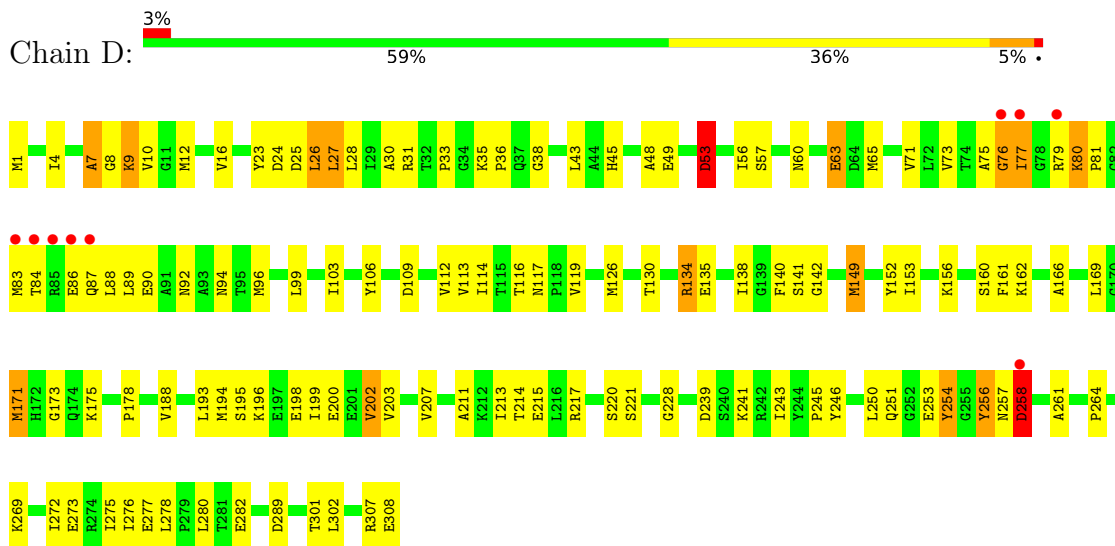
3 Residue-property plots [i](#)

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: Malate dehydrogenase

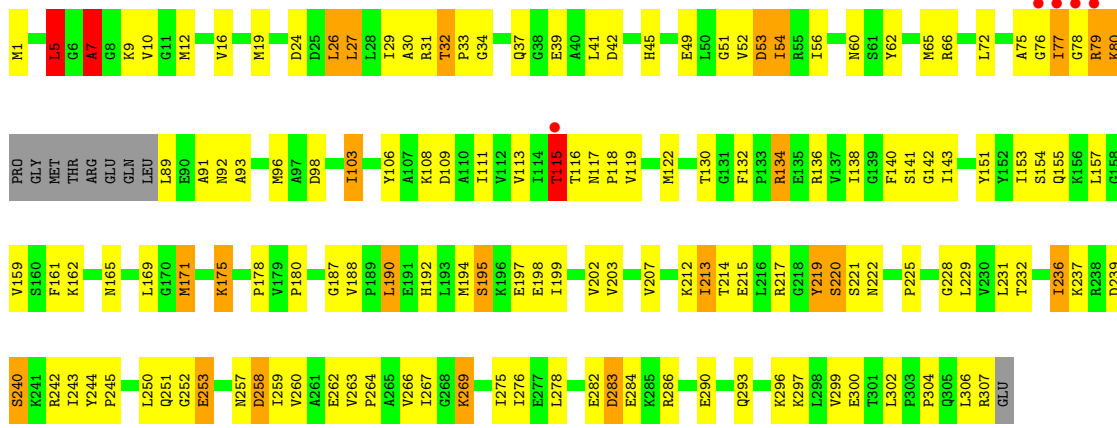


- Molecule 1: Malate dehydrogenase

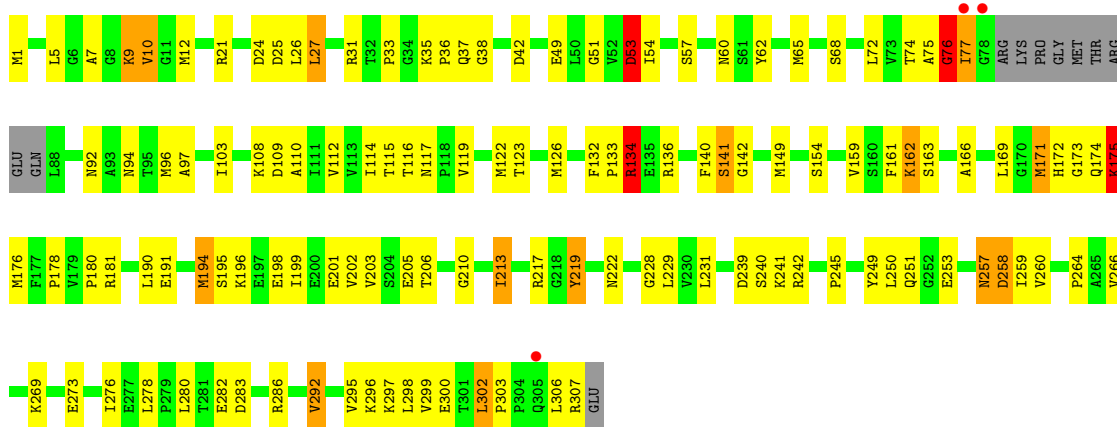


- Molecule 1: Malate dehydrogenase





• Molecule 1: Malate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.17Å 84.01Å 216.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.87 5.15 – 2.87	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.92-2.87) 99.8 (5.15-2.87)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 2.89Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.205 , 0.240 0.195 , 0.230	Depositor DCC
R_{free} test set	2707 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtrriage
Anisotropy	0.086	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.86 , 113.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9292	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.49% 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 i

5.1 Standard geometry i

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	0.65	7/2295 (0.3%)	1.08	20/3100 (0.6%)
1	B	0.88	13/2322 (0.6%)	1.31	37/3136 (1.2%)
1	C	0.70	10/2307 (0.4%)	1.04	16/3114 (0.5%)
1	D	0.63	4/2381 (0.2%)	0.97	16/3215 (0.5%)
All	All	0.72	34/9305 (0.4%)	1.10	89/12565 (0.7%)

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	3
1	B	0	3
1	C	0	4
1	D	0	2
All	All	0	12

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	254	TYR	C-N	8.37	1.48	1.33
1	C	5	LEU	C-N	7.81	1.47	1.33
1	C	253	GLU	C-N	-7.72	1.16	1.34
1	A	134	ARG	C-O	-7.71	1.08	1.23
1	B	217	ARG	C-N	-7.65	1.19	1.33
1	C	7	ALA	C-N	7.49	1.46	1.33
1	D	135	GLU	C-N	-7.40	1.17	1.34
1	B	106	TYR	C-N	7.11	1.50	1.34
1	B	240	SER	C-N	-6.81	1.18	1.34
1	C	258	ASP	C-N	6.77	1.49	1.34
1	B	52	VAL	C-N	-6.63	1.18	1.34
1	B	188	VAL	C-N	-6.60	1.21	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	195	SER	C-O	6.45	1.35	1.23
1	B	179	VAL	C-O	6.29	1.35	1.23
1	D	258	ASP	C-N	6.25	1.48	1.34
1	B	67	GLY	C-N	6.24	1.48	1.34
1	B	200	GLU	C-N	6.21	1.48	1.34
1	A	123	THR	C-N	-6.05	1.20	1.34
1	C	284	GLU	C-N	5.98	1.47	1.34
1	C	219	TYR	C-N	-5.93	1.20	1.34
1	D	38	GLY	C-N	-5.88	1.20	1.34
1	B	254	TYR	C-N	5.80	1.43	1.33
1	C	9	LYS	C-N	-5.78	1.20	1.34
1	B	119	VAL	N-CA	-5.75	1.34	1.46
1	A	257	ASN	C-N	-5.72	1.20	1.34
1	A	57	SER	C-N	5.61	1.43	1.33
1	A	292	VAL	C-N	-5.49	1.21	1.34
1	C	115	THR	C-N	5.46	1.46	1.34
1	B	197	GLU	C-N	5.34	1.46	1.34
1	A	112	VAL	C-N	-5.33	1.21	1.34
1	A	21	ARG	C-N	-5.33	1.23	1.33
1	C	54	ILE	C-N	5.25	1.46	1.34
1	B	118	PRO	C-O	-5.14	1.12	1.23
1	B	8	GLY	C-N	5.12	1.45	1.34

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	7	ALA	O-C-N	-12.69	101.64	123.20
1	C	115	THR	O-C-N	-12.28	103.06	122.70
1	C	219	TYR	O-C-N	-11.90	103.66	122.70
1	B	201	GLU	O-C-N	-11.79	103.84	122.70
1	C	240	SER	O-C-N	-11.14	104.88	122.70
1	B	224	GLY	C-N-CD	10.59	150.65	128.40
1	D	134	ARG	O-C-N	10.35	139.25	122.70
1	B	217	ARG	O-C-N	-10.32	105.66	123.20
1	B	106	TYR	O-C-N	-9.95	106.78	122.70
1	A	196	LYS	O-C-N	9.50	137.90	122.70
1	B	254	TYR	O-C-N	-9.44	107.16	123.20
1	D	135	GLU	O-C-N	-9.41	107.64	122.70
1	A	134	ARG	O-C-N	-8.46	109.17	122.70
1	B	223	TYR	CB-CG-CD2	8.32	125.99	121.00
1	D	258	ASP	O-C-N	-8.26	109.48	122.70
1	C	219	TYR	C-N-CA	8.21	142.22	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	134	ARG	CA-C-N	-8.19	99.18	117.20
1	C	258	ASP	O-C-N	-8.04	109.84	122.70
1	B	254	TYR	CA-C-N	7.66	131.52	116.20
1	C	5	LEU	C-N-CA	-7.58	106.39	122.30
1	A	163	SER	O-C-N	-7.50	110.70	122.70
1	A	38	GLY	O-C-N	-7.46	110.76	122.70
1	C	219	TYR	CA-C-N	7.43	133.55	117.20
1	C	195	SER	O-C-N	-7.38	110.89	122.70
1	C	240	SER	CA-C-N	7.33	133.32	117.20
1	B	201	GLU	CA-C-N	7.31	133.28	117.20
1	A	196	LYS	CA-C-N	-7.20	101.35	117.20
1	B	253	GLU	O-C-N	7.20	134.21	122.70
1	B	202	VAL	C-N-CA	-7.17	103.78	121.70
1	B	225	PRO	CA-N-CD	-7.12	101.53	111.50
1	B	307	ARG	O-C-N	-7.07	111.39	122.70
1	B	111	ILE	O-C-N	-7.07	111.39	122.70
1	D	7	ALA	O-C-N	-7.02	111.26	123.20
1	B	197	GLU	CA-C-N	-7.02	101.76	117.20
1	D	135	GLU	CA-C-N	6.90	132.39	117.20
1	D	77	ILE	CA-C-N	6.84	129.87	116.20
1	A	10	VAL	O-C-N	6.81	134.78	123.20
1	D	254	TYR	O-C-N	-6.69	111.82	123.20
1	B	119	VAL	O-C-N	6.68	133.38	122.70
1	B	197	GLU	O-C-N	6.68	133.38	122.70
1	A	175	LYS	O-C-N	-6.67	112.03	122.70
1	A	162	LYS	O-C-N	6.62	133.30	122.70
1	B	52	VAL	C-N-CA	6.52	137.99	121.70
1	C	98	ASP	O-C-N	-6.49	112.31	122.70
1	C	213	ILE	CA-C-N	-6.46	103.00	117.20
1	D	77	ILE	O-C-N	-6.41	112.30	123.20
1	A	10	VAL	C-N-CA	-6.36	108.94	122.30
1	D	134	ARG	C-N-CA	-6.30	105.95	121.70
1	B	180	PRO	O-C-N	-6.28	112.66	122.70
1	B	53	ASP	CB-CG-OD2	6.18	123.86	118.30
1	B	254	TYR	C-N-CA	6.04	134.99	122.30
1	A	194	MET	CA-C-N	-6.01	103.97	117.20
1	B	179	VAL	O-C-N	5.96	132.42	121.10
1	B	106	TYR	CA-C-N	5.95	130.29	117.20
1	B	223	TYR	CB-CG-CD1	-5.82	117.51	121.00
1	A	10	VAL	CA-C-N	-5.81	104.58	116.20
1	D	53	ASP	O-C-N	-5.78	113.46	122.70
1	B	253	GLU	CA-C-N	-5.73	104.60	117.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	52	VAL	O-C-N	-5.73	113.54	122.70
1	C	32	THR	O-C-N	5.69	131.92	121.10
1	C	213	ILE	C-N-CA	-5.65	107.57	121.70
1	B	119	VAL	CA-C-N	-5.64	104.79	117.20
1	A	219	TYR	O-C-N	5.64	131.72	122.70
1	B	69	ASP	CB-CG-OD2	-5.63	113.23	118.30
1	D	254	TYR	CA-C-N	5.61	127.42	116.20
1	B	118	PRO	CA-C-N	5.58	129.47	117.20
1	D	135	GLU	C-N-CA	5.56	135.60	121.70
1	C	240	SER	C-N-CA	5.52	135.50	121.70
1	A	76	GLY	CA-C-N	-5.46	105.19	117.20
1	A	133	PRO	O-C-N	5.46	131.43	122.70
1	D	141	SER	O-C-N	-5.45	113.93	123.20
1	A	162	LYS	CA-C-N	-5.40	105.31	117.20
1	B	67	GLY	C-N-CA	-5.34	108.34	121.70
1	B	107	ALA	CB-CA-C	5.34	118.11	110.10
1	A	163	SER	CA-C-N	5.34	128.95	117.20
1	A	38	GLY	CA-C-N	5.33	128.93	117.20
1	A	27	LEU	O-C-N	-5.33	114.18	122.70
1	B	111	ILE	CA-C-N	5.29	128.84	117.20
1	A	258	ASP	C-N-CA	5.25	134.83	121.70
1	B	180	PRO	C-N-CA	5.24	134.80	121.70
1	B	119	VAL	C-N-CA	-5.22	108.65	121.70
1	A	133	PRO	CA-C-N	-5.18	105.80	117.20
1	D	258	ASP	CA-C-O	5.14	130.89	120.10
1	C	115	THR	CA-C-O	5.10	130.80	120.10
1	B	8	GLY	CA-C-N	-5.09	105.99	117.20
1	B	202	VAL	CA-C-N	-5.09	106.00	117.20
1	B	204	SER	CA-C-N	-5.06	106.08	117.20
1	D	256	TYR	CB-CG-CD1	-5.03	117.98	121.00
1	B	217	ARG	CA-C-N	5.01	126.21	116.20

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	134	ARG	Sidechain
1	A	194	MET	Mainchain
1	A	76	GLY	Mainchain
1	B	201	GLU	Mainchain
1	B	202	VAL	Mainchain
1	B	278	LEU	Mainchain

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Mol	Chain	Res	Type	Group
1	C	115	THR	Mainchain
1	C	219	TYR	Mainchain
1	C	258	ASP	Mainchain
1	C	7	ALA	Mainchain
1	D	258	ASP	Mainchain
1	D	8	GLY	Mainchain

5.2 Too-close contacts [i](#)

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	2262	0	2334	111	1
1	B	2289	0	2353	118	2
1	C	2274	0	2349	146	0
1	D	2346	0	2422	132	1
2	A	28	0	0	1	0
2	B	42	0	0	1	0
2	C	17	0	0	0	0
2	D	34	0	0	1	0
All	All	9292	0	9458	475	3

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 (475) 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:B:174:GLN:NE2	1:B:175:LYS:HZ2	1.34	1.24
1:B:174:GLN:NE2	1:B:175:LYS:NZ	1.93	1.17
1:B:76:GLY:HA3	1:B:116:THR:HG23	1.32	1.09
1:A:76:GLY:HA3	1:A:116:THR:HG23	1.33	1.06
1:A:259:ILE:HD13	1:A:292:VAL:HG13	1.44	1.00
1:B:242:ARG:NH2	1:D:53:ASP:OD1	1.95	0.98
1:D:80:LYS:O	1:D:83:MET:HB2	1.63	0.96
1:B:118:PRO:O	1:B:122:MET:HG2	1.68	0.93
1:D:134:ARG:HH11	1:D:251:GLN:NE2	1.67	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:84:THR:HB	1:D:87:GLN:HB2	1.50	0.91
1:D:84:THR:HB	1:D:87:GLN:CB	2.04	0.88
1:D:243:ILE:HD13	1:D:276:ILE:CD1	2.03	0.87
1:C:232:THR:O	1:C:236:ILE:HD13	1.73	0.87
1:C:53:ASP:OD1	1:A:242:ARG:HD2	1.75	0.87
1:B:153:ILE:CD1	1:B:190:LEU:HD11	2.05	0.87
1:B:154:SER:HB2	1:B:159:VAL:O	1.76	0.86
1:A:195:SER:OG	1:A:198:GLU:HG3	1.76	0.86
1:D:26:LEU:HD21	1:D:56:ILE:HG12	1.59	0.85
1:B:153:ILE:HD13	1:B:190:LEU:HD11	1.59	0.84
1:D:134:ARG:HH11	1:D:251:GLN:HE21	1.24	0.84
1:D:140:PHE:CZ	1:D:228:GLY:HA3	2.11	0.84
1:B:77:ILE:HG22	1:B:92:ASN:OD1	1.76	0.84
1:C:296:LYS:O	1:C:300:GLU:HG3	1.79	0.83
1:B:174:GLN:HE21	1:B:175:LYS:NZ	1.67	0.82
1:C:297:LYS:HB3	1:C:297:LYS:NZ	1.95	0.82
1:C:142:GLY:HA3	1:C:245:PRO:HG2	1.62	0.81
1:B:174:GLN:HE21	1:B:175:LYS:HZ2	0.82	0.81
1:B:171:MET:HE2	1:B:173:GLY:N	1.95	0.81
1:C:143:ILE:HD11	1:C:244:TYR:HB3	1.63	0.80
1:A:75:ALA:O	1:A:76:GLY:O	1.99	0.79
1:B:12:MET:SD	1:B:39:GLU:HG2	2.22	0.79
1:A:303:PRO:HD2	1:A:306:LEU:HD12	1.62	0.79
1:A:76:GLY:HA3	1:A:116:THR:CG2	2.12	0.79
1:A:259:ILE:HD12	1:A:296:LYS:HG3	1.63	0.79
1:C:236:ILE:HD11	1:C:267:ILE:CD1	2.13	0.79
1:C:269:LYS:HB3	1:C:269:LYS:NZ	1.95	0.79
1:B:204:SER:O	1:B:208:ASN:HB2	1.82	0.79
1:C:154:SER:HB2	1:C:159:VAL:O	1.83	0.79
1:D:171:MET:HE3	1:D:173:GLY:N	1.98	0.79
1:A:9:LYS:HB2	1:A:9:LYS:NZ	1.98	0.78
1:D:213:ILE:HD11	1:D:221:SER:HB2	1.64	0.78
1:A:266:VAL:HG23	1:A:276:ILE:HD11	1.65	0.78
1:D:138:ILE:HG13	1:D:272:ILE:HD11	1.65	0.77
1:A:117:ASN:OD1	1:A:119:VAL:HG23	1.84	0.77
1:B:9:LYS:HD3	1:B:219:TYR:CE1	2.19	0.77
1:B:76:GLY:CA	1:B:116:THR:HG23	2.13	0.77
1:D:117:ASN:OD1	1:D:119:VAL:HG23	1.85	0.77
1:A:171:MET:HE1	1:A:175:LYS:H	1.50	0.76
1:C:239:ASP:HB2	1:C:269:LYS:HB2	1.66	0.76
1:C:77:ILE:HD13	1:C:78:GLY:N	2.01	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:134:ARG:HD2	1:D:251:GLN:CD	2.07	0.75
1:B:174:GLN:NE2	1:B:175:LYS:HZ1	1.85	0.74
1:D:239:ASP:OD1	1:D:241:LYS:NZ	2.15	0.73
1:C:140:PHE:CE2	1:C:225:PRO:HA	2.23	0.73
1:A:115:THR:HG21	1:A:229:LEU:HD11	1.71	0.73
1:B:213:ILE:O	1:B:217:ARG:HG2	1.89	0.73
1:C:140:PHE:CZ	1:C:228:GLY:HA3	2.24	0.73
1:D:140:PHE:CE1	1:D:228:GLY:HA3	2.24	0.73
1:B:63:GLU:O	1:B:66:ARG:HG3	1.89	0.72
1:C:169:LEU:HD11	1:C:278:LEU:HD12	1.70	0.72
1:B:53:ASP:OD2	1:A:162:LYS:HG3	1.89	0.71
1:D:153:ILE:HG23	1:D:194:MET:CE	2.20	0.71
1:A:171:MET:CE	1:A:175:LYS:H	2.02	0.71
1:D:80:LYS:HG2	1:D:81:PRO:HD2	1.73	0.71
1:A:1:MET:HA	1:A:24:ASP:OD2	1.90	0.71
1:C:103:ILE:HD11	1:C:130:THR:HG22	1.71	0.71
1:A:154:SER:HB2	1:A:159:VAL:O	1.90	0.71
1:C:7:ALA:HA	1:C:12:MET:HE3	1.74	0.70
1:C:141:SER:HB2	1:C:262:GLU:OE1	1.92	0.69
1:A:7:ALA:HA	1:A:12:MET:HE3	1.73	0.69
1:A:249:TYR:CZ	1:A:258:ASP:HA	2.29	0.68
1:C:77:ILE:HG22	1:C:92:ASN:OD1	1.93	0.68
1:D:79:ARG:HE	1:D:83:MET:HB3	1.59	0.68
1:A:115:THR:CG2	1:A:229:LEU:HD11	2.24	0.68
1:C:140:PHE:CE1	1:C:228:GLY:HA3	2.29	0.68
1:C:79:ARG:HH11	1:C:79:ARG:HB2	1.58	0.68
1:A:210:GLY:HA2	1:A:213:ILE:HD11	1.76	0.67
1:A:286:ARG:HB3	1:A:286:ARG:NH1	2.09	0.67
1:B:304:PRO:HA	1:B:307:ARG:HG2	1.76	0.67
1:B:307:ARG:O	1:B:308:GLU:HB2	1.95	0.67
1:D:152:TYR:HB3	1:D:202:VAL:HG12	1.77	0.66
1:C:213:ILE:HD11	1:C:221:SER:HB2	1.76	0.66
1:D:195:SER:OG	1:D:198:GLU:HG3	1.96	0.66
1:A:282:GLU:HG3	1:A:283:ASP:N	2.11	0.66
1:C:269:LYS:HB3	1:C:269:LYS:HZ3	1.61	0.66
1:A:178:PRO:HG3	1:A:203:VAL:HG22	1.77	0.66
1:C:199:ILE:O	1:C:203:VAL:HG23	1.96	0.65
1:A:199:ILE:O	1:A:203:VAL:HG23	1.96	0.65
1:B:242:ARG:CZ	1:D:53:ASP:OD1	2.43	0.65
1:B:174:GLN:HE22	1:B:175:LYS:NZ	1.91	0.65
1:B:171:MET:HE2	1:B:173:GLY:CA	2.26	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:196:LYS:O	1:D:200:GLU:HG3	1.96	0.65
1:A:171:MET:HE3	1:A:173:GLY:N	2.11	0.65
1:C:297:LYS:HB3	1:C:297:LYS:HZ3	1.62	0.65
1:A:240:SER:OG	1:A:242:ARG:HD3	1.96	0.65
1:D:89:LEU:HG	1:D:302:LEU:HD11	1.78	0.65
1:C:293:GLN:O	1:C:296:LYS:HB3	1.97	0.65
1:B:162:LYS:HD3	1:A:53:ASP:OD1	1.96	0.65
1:C:79:ARG:HG2	1:C:214:THR:HG21	1.78	0.64
1:B:254:TYR:N	1:B:277:GLU:OE1	2.28	0.64
1:D:71:VAL:HG11	1:D:103:ILE:CD1	2.27	0.64
1:C:169:LEU:HD21	1:C:264:PRO:HD3	1.80	0.64
1:D:253:GLU:HG2	1:D:275:ILE:HB	1.80	0.64
1:B:181:ARG:NH1	1:B:182:LEU:HD21	2.12	0.64
1:D:269:LYS:NZ	1:D:269:LYS:HB3	2.13	0.64
1:B:217:ARG:HH22	1:B:222:ASN:N	1.95	0.64
1:B:181:ARG:O	1:B:182:LEU:HD23	1.98	0.63
1:D:7:ALA:HA	1:D:12:MET:HE3	1.78	0.63
1:A:134:ARG:HH12	1:A:253:GLU:CD	2.01	0.63
1:D:243:ILE:HD13	1:D:276:ILE:HD13	1.80	0.63
1:A:134:ARG:NH1	1:A:253:GLU:OE1	2.31	0.63
1:D:162:LYS:HG3	1:C:53:ASP:OD2	1.97	0.63
1:D:171:MET:HE3	1:D:173:GLY:H	1.63	0.62
1:A:259:ILE:CD1	1:A:292:VAL:HG13	2.25	0.62
1:A:282:GLU:HG3	1:A:283:ASP:H	1.64	0.62
1:B:167:ILE:HD11	1:C:187:GLY:HA2	1.80	0.62
1:A:72:LEU:HD22	1:A:229:LEU:HD22	1.82	0.62
1:D:84:THR:HB	1:D:87:GLN:HB3	1.82	0.62
1:A:269:LYS:NZ	1:A:269:LYS:HB3	2.14	0.62
1:C:7:ALA:HA	1:C:12:MET:CE	2.29	0.62
1:B:231:LEU:HD11	1:A:49:GLU:O	2.00	0.62
1:D:241:LYS:NZ	1:D:273:GLU:OE2	2.31	0.62
1:B:213:ILE:HD11	1:B:221:SER:HB2	1.80	0.62
1:D:241:LYS:CD	1:D:273:GLU:OE2	2.47	0.62
1:B:188:VAL:HG22	1:C:278:LEU:HD23	1.82	0.61
1:D:71:VAL:HG11	1:D:103:ILE:HD11	1.82	0.61
1:D:169:LEU:HD11	1:D:278:LEU:HD12	1.82	0.61
1:A:9:LYS:HB2	1:A:9:LYS:HZ2	1.65	0.61
1:C:118:PRO:O	1:C:122:MET:HG2	2.00	0.61
1:B:135:GLU:O	1:B:271:GLY:HA3	2.00	0.61
1:D:138:ILE:CG1	1:D:272:ILE:HD11	2.30	0.61
1:D:250:LEU:HD11	1:D:261:ALA:HB3	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:LYS:HD3	1:A:219:TYR:CE1	2.36	0.61
1:B:140:PHE:CZ	1:B:228:GLY:HA3	2.35	0.61
1:B:162:LYS:CD	1:A:53:ASP:OD1	2.49	0.60
1:B:217:ARG:NH2	1:B:221:SER:HA	2.16	0.60
1:B:92:ASN:C	1:B:122:MET:HE1	2.22	0.60
1:D:153:ILE:HG23	1:D:194:MET:HE1	1.83	0.60
1:A:134:ARG:HD2	1:A:251:GLN:NE2	2.15	0.60
1:A:142:GLY:HA3	1:A:245:PRO:HG2	1.83	0.60
1:C:236:ILE:HD11	1:C:267:ILE:HD11	1.84	0.60
1:C:236:ILE:HD11	1:C:267:ILE:HD13	1.82	0.60
1:B:49:GLU:O	1:A:231:LEU:HD11	2.02	0.60
1:C:72:LEU:HD23	1:C:113:VAL:HG22	1.83	0.60
1:D:24:ASP:OD1	1:D:25:ASP:N	2.33	0.59
1:D:171:MET:HE1	1:D:175:LYS:H	1.66	0.59
1:B:242:ARG:NE	1:D:53:ASP:OD1	2.34	0.59
1:C:92:ASN:O	1:C:96:MET:HG2	2.02	0.59
1:C:66:ARG:HG3	1:C:66:ARG:HH11	1.67	0.59
1:D:199:ILE:O	1:D:203:VAL:HG23	2.03	0.59
1:C:171:MET:HE1	1:C:175:LYS:H	1.68	0.59
1:C:293:GLN:OE1	1:C:296:LYS:HD2	2.03	0.59
1:D:171:MET:HE1	1:D:175:LYS:N	2.18	0.59
1:B:217:ARG:HH22	1:B:222:ASN:H	1.50	0.58
1:B:92:ASN:CB	1:B:122:MET:HE1	2.32	0.58
1:A:276:ILE:HD12	1:A:276:ILE:N	2.17	0.58
1:A:201:GLU:O	1:A:205:GLU:HG2	2.03	0.58
1:A:140:PHE:CZ	1:A:228:GLY:HA3	2.39	0.58
1:D:7:ALA:HB3	1:D:30:ALA:HB2	1.84	0.58
1:C:195:SER:OG	1:C:198:GLU:HG3	2.04	0.58
1:D:301:THR:HG22	1:D:301:THR:O	2.03	0.58
1:C:194:MET:HB2	1:C:199:ILE:HD11	1.86	0.57
1:B:281:THR:O	1:B:285:LYS:HD3	2.05	0.57
1:C:213:ILE:O	1:C:217:ARG:HG2	2.03	0.57
1:A:296:LYS:O	1:A:300:GLU:HG3	2.04	0.57
1:B:142:GLY:HA3	1:B:245:PRO:HG2	1.85	0.57
1:B:201:GLU:HB3	2:B:338:HOH:O	2.05	0.57
1:B:86:GLU:OE2	1:B:301:THR:HG21	2.05	0.57
1:C:259:ILE:HG13	1:C:260:VAL:H	1.70	0.57
1:C:297:LYS:NZ	1:C:297:LYS:CB	2.66	0.57
1:B:92:ASN:HB3	1:B:122:MET:HE1	1.85	0.56
1:D:217:ARG:CD	1:C:39:GLU:HG3	2.34	0.56
1:D:302:LEU:O	1:D:307:ARG:NH1	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:ARG:HB3	1:C:220:SER:HB3	1.87	0.56
1:A:68:SER:O	1:A:110:ALA:HB2	2.05	0.56
1:C:199:ILE:O	1:C:202:VAL:HG22	2.04	0.56
1:D:134:ARG:NH1	1:D:251:GLN:HE21	1.98	0.56
1:D:156:LYS:HE3	1:D:198:GLU:OE1	2.06	0.56
1:C:79:ARG:HG3	1:C:80:LYS:H	1.71	0.56
1:D:138:ILE:N	1:D:138:ILE:HD12	2.21	0.55
1:C:194:MET:HB3	1:C:198:GLU:HB2	1.87	0.55
1:C:250:LEU:O	1:C:257:ASN:HA	2.06	0.55
1:B:222:ASN:C	1:B:225:PRO:HD2	2.26	0.55
1:B:162:LYS:HD2	1:A:51:GLY:O	2.05	0.55
1:D:89:LEU:HG	1:D:302:LEU:CD1	2.36	0.55
1:B:217:ARG:NH2	1:B:222:ASN:OD1	2.40	0.55
1:D:213:ILE:HG22	1:C:41:LEU:HD23	1.89	0.55
1:D:26:LEU:CD2	1:D:56:ILE:HG12	2.34	0.55
1:D:188:VAL:HB	1:D:193:LEU:HD11	1.87	0.55
1:A:202:VAL:HG23	1:A:203:VAL:N	2.22	0.54
1:C:259:ILE:HG13	1:C:260:VAL:N	2.23	0.54
1:A:286:ARG:HB3	1:A:286:ARG:HH11	1.72	0.54
1:A:92:ASN:O	1:A:96:MET:HG2	2.07	0.54
1:C:169:LEU:CD2	1:C:264:PRO:HD3	2.36	0.54
1:A:33:PRO:HA	1:A:60:ASN:OD1	2.06	0.54
1:D:80:LYS:HD2	1:D:81:PRO:O	2.07	0.54
1:B:222:ASN:O	1:B:225:PRO:HD2	2.07	0.54
1:D:134:ARG:HD2	1:D:251:GLN:NE2	2.22	0.54
1:D:307:ARG:O	1:D:308:GLU:HB2	2.07	0.54
1:A:114:ILE:HD11	1:A:126:MET:HG2	1.90	0.54
1:A:171:MET:HE1	1:A:174:GLN:N	2.22	0.54
1:C:153:ILE:HD13	1:C:202:VAL:HG11	1.88	0.53
1:C:253:GLU:HG2	1:C:275:ILE:HB	1.91	0.53
1:B:169:LEU:HD11	1:B:278:LEU:HD12	1.90	0.53
1:B:7:ALA:HA	1:B:12:MET:CE	2.39	0.53
1:B:77:ILE:HG22	1:B:92:ASN:CG	2.28	0.53
1:C:66:ARG:HG3	1:C:66:ARG:NH1	2.23	0.53
1:C:108:LYS:HA	1:C:132:PHE:CZ	2.44	0.53
1:C:293:GLN:NE2	1:C:296:LYS:HD2	2.22	0.53
1:C:7:ALA:HB3	1:C:30:ALA:HB2	1.91	0.53
1:A:77:ILE:O	1:A:77:ILE:CG2	2.57	0.53
1:D:162:LYS:HD3	1:C:51:GLY:O	2.09	0.53
1:A:302:LEU:O	1:A:307:ARG:NH1	2.42	0.53
1:D:149:MET:HG2	1:D:166:ALA:HB1	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:ARG:HH11	1:C:79:ARG:CB	2.20	0.52
1:C:142:GLY:CA	1:C:245:PRO:HG2	2.37	0.52
1:C:171:MET:CE	1:C:175:LYS:H	2.22	0.52
1:B:62:TYR:O	1:B:65:MET:HG2	2.09	0.52
1:B:217:ARG:NH2	1:B:222:ASN:N	2.58	0.52
1:B:53:ASP:OD2	1:A:162:LYS:CG	2.56	0.52
1:C:115:THR:HG23	1:C:225:PRO:HB3	1.91	0.52
1:B:7:ALA:HA	1:B:12:MET:HE3	1.91	0.52
1:B:116:THR:O	1:B:119:VAL:HG12	2.10	0.52
1:D:90:GLU:O	1:D:94:ASN:ND2	2.43	0.52
1:A:171:MET:HE3	1:A:173:GLY:H	1.74	0.52
1:D:65:MET:SD	1:D:103:ILE:HD13	2.49	0.52
1:D:77:ILE:HG22	1:D:92:ASN:OD1	2.09	0.52
1:B:171:MET:CE	1:B:173:GLY:HA3	2.39	0.52
1:B:299:VAL:O	1:B:302:LEU:HB2	2.10	0.52
1:D:307:ARG:O	1:D:308:GLU:CB	2.57	0.52
1:A:5:LEU:HD21	1:A:103:ILE:HD11	1.92	0.52
1:B:88:LEU:HG	1:B:92:ASN:ND2	2.25	0.51
1:B:153:ILE:HD11	1:B:190:LEU:HD11	1.87	0.51
1:C:282:GLU:HG3	1:C:283:ASP:N	2.25	0.51
1:B:167:ILE:CD1	1:C:187:GLY:HA2	2.41	0.51
1:D:243:ILE:CD1	1:D:276:ILE:CD1	2.84	0.51
1:C:62:TYR:O	1:C:65:MET:HG2	2.10	0.51
1:A:213:ILE:O	1:A:217:ARG:HG2	2.11	0.51
1:C:1:MET:HA	1:C:24:ASP:OD2	2.10	0.51
1:C:26:LEU:HD22	1:C:54:ILE:HD12	1.93	0.51
1:D:4:ILE:HB	1:D:28:LEU:HD13	1.93	0.51
1:C:142:GLY:HA3	1:C:245:PRO:CG	2.39	0.51
1:D:1:MET:HA	1:D:24:ASP:OD2	2.09	0.51
1:C:293:GLN:CD	1:C:296:LYS:HD2	2.32	0.50
1:D:134:ARG:CD	1:D:251:GLN:NE2	2.74	0.50
1:D:217:ARG:HD3	1:C:39:GLU:HG3	1.93	0.50
1:D:75:ALA:O	1:D:76:GLY:O	2.29	0.50
1:D:199:ILE:O	1:D:202:VAL:HG23	2.11	0.50
1:D:217:ARG:HD2	1:C:39:GLU:HG3	1.92	0.50
1:A:134:ARG:HD2	1:A:251:GLN:CD	2.32	0.50
1:D:243:ILE:CD1	1:D:276:ILE:HD11	2.42	0.50
1:D:257:ASN:O	1:D:258:ASP:C	2.49	0.50
1:C:10:VAL:HA	1:C:222:ASN:HB2	1.93	0.50
1:B:269:LYS:NZ	1:B:269:LYS:HB3	2.27	0.50
1:C:92:ASN:HB3	1:C:116:THR:HG21	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:LEU:HD11	1:B:302:LEU:HD11	1.94	0.50
1:C:89:LEU:C	1:C:91:ALA:H	2.15	0.50
1:C:190:LEU:HD13	1:C:199:ILE:HD12	1.93	0.50
1:A:9:LYS:HB2	1:A:9:LYS:HZ3	1.72	0.50
1:B:96:MET:HG2	1:B:122:MET:CE	2.42	0.49
1:D:113:VAL:HA	1:D:138:ILE:O	2.12	0.49
1:C:198:GLU:O	1:C:202:VAL:HG13	2.12	0.49
1:D:254:TYR:CZ	1:D:277:GLU:HA	2.47	0.49
1:C:203:VAL:O	1:C:207:VAL:HG23	2.12	0.49
1:C:286:ARG:NH1	1:C:290:GLU:OE2	2.44	0.49
1:D:7:ALA:HA	1:D:12:MET:CE	2.42	0.49
1:A:149:MET:CE	1:A:166:ALA:HB1	2.43	0.49
1:B:93:ALA:HA	1:B:122:MET:HE3	1.94	0.49
1:D:76:GLY:HA3	1:D:116:THR:HG23	1.94	0.49
1:B:171:MET:CE	1:B:173:GLY:N	2.71	0.49
1:C:194:MET:HB2	1:C:199:ILE:CD1	2.43	0.49
1:D:195:SER:O	1:D:199:ILE:HG13	2.12	0.49
1:A:117:ASN:HA	1:A:119:VAL:N	2.28	0.49
1:B:278:LEU:HD23	1:C:188:VAL:HG22	1.95	0.49
1:B:171:MET:CE	1:B:173:GLY:CA	2.91	0.49
1:B:194:MET:HG2	1:B:198:GLU:OE1	2.13	0.49
1:D:160:SER:HB2	1:A:240:SER:O	2.12	0.49
1:D:27:LEU:HD22	1:D:28:LEU:N	2.27	0.48
1:D:117:ASN:HA	1:D:119:VAL:N	2.28	0.48
1:C:34:GLY:H	1:C:60:ASN:HD21	1.59	0.48
1:C:113:VAL:HA	1:C:138:ILE:O	2.13	0.48
1:B:77:ILE:HG13	1:B:78:GLY:N	2.28	0.48
1:D:48:ALA:HA	1:C:161:PHE:CE1	2.48	0.48
1:B:48:ALA:HA	1:A:161:PHE:CE1	2.48	0.48
1:D:92:ASN:O	1:D:96:MET:HG2	2.14	0.48
1:D:213:ILE:CG2	1:C:41:LEU:HD23	2.42	0.48
1:C:269:LYS:HB3	1:C:269:LYS:HZ2	1.76	0.48
1:A:241:LYS:NZ	1:A:273:GLU:OE2	2.35	0.48
1:B:41:LEU:HD23	1:A:213:ILE:HG23	1.95	0.48
1:D:214:THR:OG1	1:D:220:SER:HA	2.14	0.48
1:C:26:LEU:CD2	1:C:54:ILE:HD12	2.43	0.48
1:B:168:VAL:HA	1:B:178:PRO:HA	1.95	0.48
1:B:181:ARG:NH2	1:B:279:PRO:O	2.47	0.48
1:D:79:ARG:NH2	1:D:83:MET:O	2.45	0.48
1:D:153:ILE:HG23	1:D:194:MET:HE2	1.96	0.48
1:D:241:LYS:CE	1:D:273:GLU:OE2	2.62	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:ARG:NH2	1:C:106:TYR:HD1	2.12	0.48
1:C:79:ARG:HG3	1:C:80:LYS:N	2.28	0.48
1:C:27:LEU:HD22	1:C:29:ILE:HG13	1.96	0.48
1:D:7:ALA:HB3	1:D:30:ALA:CB	2.43	0.47
1:D:35:LYS:HB3	1:D:36:PRO:HD3	1.95	0.47
1:B:269:LYS:NZ	1:B:269:LYS:CB	2.77	0.47
1:D:142:GLY:HA3	1:D:245:PRO:HG2	1.96	0.47
1:D:169:LEU:CD2	1:D:264:PRO:HD3	2.44	0.47
1:C:243:ILE:CD1	1:C:266:VAL:HG22	2.44	0.47
1:D:63:GLU:HG3	1:D:106:TYR:CZ	2.48	0.47
1:C:153:ILE:HD11	1:C:202:VAL:HG21	1.97	0.47
1:C:134:ARG:HH22	1:C:253:GLU:CD	2.18	0.47
1:B:167:ILE:HD12	1:B:182:LEU:CB	2.45	0.47
1:C:109:ASP:HA	1:C:136:ARG:NH1	2.30	0.47
1:D:269:LYS:NZ	1:D:269:LYS:CB	2.78	0.47
1:C:5:LEU:HD12	1:C:29:ILE:HB	1.97	0.47
1:B:307:ARG:O	1:B:308:GLU:CB	2.61	0.47
1:C:245:PRO:HA	1:C:263:VAL:O	2.15	0.47
1:A:24:ASP:OD1	1:A:25:ASP:N	2.42	0.47
1:D:213:ILE:HD12	1:C:42:ASP:HB2	1.96	0.47
1:D:256:TYR:OH	1:D:289:ASP:OD1	2.28	0.47
1:B:96:MET:HG2	1:B:122:MET:HE3	1.97	0.46
1:A:171:MET:CE	1:A:175:LYS:N	2.75	0.46
1:C:111:ILE:HD12	1:C:111:ILE:N	2.30	0.46
1:A:5:LEU:CD2	1:A:103:ILE:HD11	2.46	0.46
1:C:117:ASN:HA	1:C:119:VAL:N	2.30	0.46
1:C:297:LYS:CB	1:C:297:LYS:HZ2	2.26	0.46
1:D:196:LYS:HE3	1:D:200:GLU:OE2	2.15	0.46
1:C:113:VAL:HG12	1:C:138:ILE:HB	1.97	0.46
1:B:175:LYS:HE2	1:B:294:ALA:HB2	1.98	0.46
1:A:269:LYS:HB3	1:A:269:LYS:HZ3	1.81	0.46
1:B:181:ARG:HH12	1:B:182:LEU:HD21	1.79	0.46
1:B:188:VAL:HG11	1:B:193:LEU:HD21	1.97	0.46
1:A:140:PHE:CE1	1:A:228:GLY:HA3	2.51	0.46
1:B:174:GLN:HE22	1:B:175:LYS:HZ1	1.56	0.46
1:D:35:LYS:HB3	1:D:36:PRO:CD	2.45	0.46
1:D:45:HIS:O	1:D:49:GLU:HG3	2.15	0.46
1:C:7:ALA:HB3	1:C:30:ALA:CB	2.46	0.46
1:C:239:ASP:CB	1:C:269:LYS:HB2	2.43	0.46
1:A:75:ALA:C	1:A:76:GLY:O	2.54	0.46
1:B:115:THR:HG23	1:B:225:PRO:HB3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:MET:HE1	1:B:173:GLY:HA3	1.98	0.45
1:B:269:LYS:HB3	1:B:269:LYS:HZ3	1.81	0.45
1:D:71:VAL:HG11	1:D:103:ILE:HD12	1.97	0.45
1:A:9:LYS:CD	1:A:219:TYR:CE1	2.99	0.45
1:B:140:PHE:CE1	1:B:228:GLY:HA3	2.51	0.45
1:C:269:LYS:NZ	1:C:269:LYS:CB	2.73	0.45
1:A:171:MET:O	1:A:176:MET:HB2	2.15	0.45
1:C:243:ILE:HG13	1:C:276:ILE:HD11	1.98	0.45
1:C:299:VAL:O	1:C:302:LEU:HB2	2.16	0.45
1:A:142:GLY:C	1:A:245:PRO:HG2	2.37	0.45
1:D:211:ALA:O	1:D:215:GLU:HG2	2.16	0.45
1:C:12:MET:O	1:C:16:VAL:HG23	2.16	0.45
1:A:26:LEU:CD2	1:A:54:ILE:HD12	2.46	0.45
1:A:141:SER:HB2	1:A:172:HIS:HB2	1.99	0.45
1:D:79:ARG:HD2	1:D:79:ARG:HA	1.83	0.45
1:C:79:ARG:CB	1:C:220:SER:HB3	2.47	0.45
1:A:191:GLU:HG3	2:A:319:HOH:O	2.16	0.45
1:D:140:PHE:CE1	1:D:228:GLY:CA	2.97	0.45
1:B:240:SER:OG	1:B:242:ARG:HG3	2.17	0.45
1:D:10:VAL:HG12	2:D:317:HOH:O	2.16	0.45
1:D:1:MET:SD	1:D:27:LEU:HB2	2.57	0.45
1:D:198:GLU:O	1:D:202:VAL:HG22	2.16	0.45
1:A:94:ASN:O	1:A:97:ALA:HB3	2.17	0.45
1:B:3:THR:HG22	1:B:5:LEU:HD13	1.99	0.45
1:D:80:LYS:O	1:D:83:MET:CB	2.52	0.45
1:C:293:GLN:HE22	1:C:296:LYS:HD2	1.82	0.45
1:A:142:GLY:CA	1:A:245:PRO:HG2	2.46	0.45
1:A:202:VAL:CG2	1:A:203:VAL:N	2.80	0.45
1:C:115:THR:OG1	1:C:229:LEU:HD11	2.16	0.44
1:C:171:MET:CE	1:C:175:LYS:N	2.80	0.44
1:B:213:ILE:CD1	1:B:221:SER:HB2	2.46	0.44
1:A:75:ALA:O	1:A:76:GLY:C	2.54	0.44
1:A:259:ILE:HG12	1:A:260:VAL:N	2.32	0.44
1:D:86:GLU:OE2	1:D:301:THR:HG21	2.17	0.44
1:D:12:MET:O	1:D:16:VAL:HG23	2.17	0.44
1:C:299:VAL:O	1:C:307:ARG:HD3	2.18	0.44
1:A:169:LEU:CD2	1:A:264:PRO:HD3	2.48	0.44
1:A:278:LEU:O	1:A:280:LEU:HD13	2.18	0.44
1:B:188:VAL:HA	1:B:189:PRO:HD3	1.76	0.44
1:B:241:LYS:HD2	1:B:273:GLU:OE1	2.18	0.44
1:D:79:ARG:HG2	1:D:88:LEU:HD22	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:203:VAL:O	1:D:207:VAL:HG23	2.18	0.44
1:C:253:GLU:CD	1:C:253:GLU:H	2.21	0.44
1:C:19:MET:O	1:C:52:VAL:HG11	2.18	0.44
1:B:242:ARG:HE	1:C:162:LYS:HD3	1.83	0.44
1:C:26:LEU:CD2	1:C:56:ILE:HG12	2.48	0.44
1:C:178:PRO:O	1:C:180:PRO:HD3	2.17	0.44
1:C:304:PRO:HA	1:C:307:ARG:HG3	1.98	0.44
1:B:66:ARG:NH1	1:B:66:ARG:HG2	2.32	0.43
1:D:71:VAL:HB	1:D:112:VAL:HG12	2.00	0.43
1:C:178:PRO:HG3	1:C:203:VAL:HG22	2.00	0.43
1:A:9:LYS:HG2	1:A:222:ASN:ND2	2.33	0.43
1:A:65:MET:SD	1:A:103:ILE:HD13	2.58	0.43
1:B:115:THR:CG2	1:B:225:PRO:HB3	2.47	0.43
1:D:171:MET:HE1	1:D:173:GLY:C	2.38	0.43
1:A:149:MET:HE2	1:A:166:ALA:HB1	2.00	0.43
1:A:36:PRO:HG2	1:A:37:GLN:OE1	2.18	0.43
1:D:188:VAL:HG22	1:A:278:LEU:HD23	2.01	0.43
1:C:26:LEU:HD23	1:C:26:LEU:H	1.82	0.43
1:C:37:GLN:HE22	1:C:60:ASN:ND2	2.16	0.43
1:C:103:ILE:HD11	1:C:130:THR:CG2	2.44	0.43
1:A:1:MET:CA	1:A:24:ASP:OD2	2.64	0.43
1:A:199:ILE:O	1:A:202:VAL:HG22	2.19	0.43
1:C:151:TYR:O	1:C:155:GLN:HG2	2.19	0.43
1:A:250:LEU:O	1:A:257:ASN:HA	2.19	0.43
1:B:96:MET:CG	1:B:122:MET:HE3	2.48	0.43
1:B:241:LYS:HE2	1:B:273:GLU:OE2	2.19	0.43
1:C:45:HIS:O	1:C:49:GLU:HG3	2.19	0.43
1:A:239:ASP:HB2	1:A:269:LYS:HB2	2.00	0.43
1:B:162:LYS:HD2	1:A:53:ASP:OD1	2.18	0.42
1:D:109:ASP:OD2	1:D:109:ASP:N	2.52	0.42
1:D:134:ARG:NH2	1:D:275:ILE:HD13	2.34	0.42
1:C:157:LEU:HG	1:C:194:MET:HE3	2.00	0.42
1:A:176:MET:CE	1:A:206:THR:HG22	2.49	0.42
1:A:286:ARG:CZ	1:A:286:ARG:CB	2.97	0.42
1:B:223:TYR:O	1:B:224:GLY:C	2.57	0.42
1:D:79:ARG:HD2	1:D:83:MET:HE2	2.00	0.42
1:A:295:VAL:O	1:A:299:VAL:HG23	2.20	0.42
1:D:308:GLU:HA	1:D:308:GLU:OE1	2.19	0.42
1:C:169:LEU:HD11	1:C:278:LEU:CD1	2.46	0.42
1:A:10:VAL:HG13	1:A:74:THR:HB	2.01	0.42
1:A:178:PRO:O	1:A:180:PRO:HD3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:84:THR:O	1:D:87:GLN:N	2.53	0.42
1:B:77:ILE:HG23	1:B:88:LEU:HD12	2.00	0.42
1:B:92:ASN:C	1:B:122:MET:CE	2.87	0.42
1:D:23:TYR:O	1:D:24:ASP:HB2	2.19	0.42
1:C:240:SER:OG	1:C:242:ARG:HG3	2.19	0.42
1:D:73:VAL:HB	1:D:114:ILE:HD13	2.02	0.42
1:D:269:LYS:HB3	1:D:269:LYS:HZ3	1.81	0.42
1:C:117:ASN:HA	1:C:118:PRO:C	2.39	0.42
1:C:243:ILE:HD11	1:C:266:VAL:HG22	2.01	0.42
1:C:297:LYS:HB3	1:C:297:LYS:HZ2	1.75	0.42
1:D:49:GLU:O	1:C:231:LEU:HD11	2.20	0.42
1:A:171:MET:HE1	1:A:173:GLY:C	2.40	0.42
1:A:169:LEU:HD11	1:A:278:LEU:HD12	2.02	0.42
1:D:27:LEU:HD23	1:D:57:SER:HB2	2.01	0.41
1:D:43:LEU:HD23	1:D:56:ILE:HG21	2.01	0.41
1:B:128:LYS:HD2	1:B:308:GLU:HG2	2.00	0.41
1:C:212:LYS:O	1:C:215:GLU:HB3	2.20	0.41
1:C:242:ARG:O	1:C:266:VAL:HA	2.20	0.41
1:B:249:TYR:OH	1:B:258:ASP:HA	2.19	0.41
1:D:33:PRO:HA	1:D:60:ASN:HD21	1.86	0.41
1:C:76:GLY:HA3	1:C:116:THR:HG23	2.01	0.41
1:A:96:MET:HG3	1:A:122:MET:SD	2.60	0.41
1:C:32:THR:HA	1:C:33:PRO:HD3	1.94	0.41
1:D:126:MET:O	1:D:130:THR:HG23	2.20	0.41
1:D:138:ILE:CG2	1:D:246:TYR:CD1	3.03	0.41
1:C:236:ILE:CD1	1:C:236:ILE:N	2.82	0.41
1:A:108:LYS:O	1:A:136:ARG:NH1	2.49	0.41
1:B:66:ARG:CG	1:B:66:ARG:HH11	2.33	0.41
1:D:9:LYS:HZ3	1:D:9:LYS:HB2	1.86	0.41
1:C:134:ARG:HH21	1:C:251:GLN:N	2.19	0.41
1:C:252:GLY:N	1:C:257:ASN:HB3	2.35	0.41
1:A:35:LYS:HB3	1:A:36:PRO:CD	2.50	0.41
1:A:297:LYS:HB3	1:A:297:LYS:HE2	1.71	0.41
1:B:9:LYS:NZ	1:B:9:LYS:HB2	2.35	0.41
1:B:96:MET:CG	1:B:122:MET:CE	2.99	0.41
1:D:80:LYS:HD2	1:D:80:LYS:C	2.41	0.41
1:C:93:ALA:HB1	1:C:306:LEU:HD13	2.03	0.41
1:B:88:LEU:HG	1:B:92:ASN:HD22	1.85	0.41
1:B:126:MET:O	1:B:130:THR:HG23	2.21	0.41
1:B:217:ARG:NH1	1:A:42:ASP:OD2	2.54	0.41
1:D:9:LYS:HB2	1:D:9:LYS:NZ	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:252:GLY:H	1:C:257:ASN:HB3	1.86	0.41
1:C:140:PHE:HZ	1:C:228:GLY:HA3	1.83	0.40
1:A:108:LYS:HA	1:A:132:PHE:CZ	2.55	0.40
1:C:72:LEU:HD23	1:C:113:VAL:CG2	2.51	0.40
1:B:7:ALA:HB3	1:B:30:ALA:HB2	2.02	0.40
1:C:52:VAL:HG12	1:C:54:ILE:HG23	2.03	0.40
1:C:192:HIS:O	1:C:192:HIS:CG	2.75	0.40
1:B:304:PRO:O	1:B:305:GLN:C	2.59	0.40
1:C:171:MET:HE2	1:C:175:LYS:CA	2.52	0.40
1:B:177:PHE:HA	1:B:178:PRO:HD2	1.91	0.40
1:B:249:TYR:CZ	1:B:258:ASP:HA	2.56	0.40
1:B:278:LEU:HB2	1:B:280:LEU:CD1	2.51	0.40
1:B:301:THR:HG22	1:B:301:THR:O	2.22	0.40
1:A:62:TYR:O	1:A:65:MET:HG2	2.21	0.40
1:A:171:MET:HE1	1:A:175:LYS:N	2.27	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:LYS:NZ	1:B:200:GLU:OE1[4_557]	1.86	0.34
1:B:133:PRO:CA	1:B:215:GLU:OE1[4_557]	1.99	0.21
1:D:87:GLN:NE2	1:A:181:ARG:NH1[3_756]	2.02	0.18

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	294/308 (96%)	276 (94%)	15 (5%)	3 (1%)	15 42
1	B	297/308 (96%)	280 (94%)	16 (5%)	1 (0%)	41 70
1	C	295/308 (96%)	279 (95%)	13 (4%)	3 (1%)	15 42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	306/308 (99%)	287 (94%)	17 (6%)	2 (1%)	22	52
All	All	1192/1232 (97%)	1122 (94%)	61 (5%)	9 (1%)	19	48

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	53	ASP
1	C	220	SER
1	B	181	ARG
1	C	53	ASP
1	C	75	ALA
1	A	53	ASP
1	A	76	GLY
1	A	141	SER
1	D	76	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	241/250 (96%)	229 (95%)	12 (5%)	24	54
1	B	244/250 (98%)	234 (96%)	10 (4%)	30	62
1	C	242/250 (97%)	224 (93%)	18 (7%)	13	36
1	D	250/250 (100%)	236 (94%)	14 (6%)	21	49
All	All	977/1000 (98%)	923 (94%)	54 (6%)	21	50

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

Mol	Chain	Res	Type
1	B	5	LEU
1	B	9	LYS
1	B	27	LEU
1	B	31	ARG
1	B	35	LYS

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Mol	Chain	Res	Type
1	B	39	GLU
1	B	66	ARG
1	B	161	PHE
1	B	165	ASN
1	B	277	GLU
1	D	9	LYS
1	D	26	LEU
1	D	27	LEU
1	D	31	ARG
1	D	63	GLU
1	D	80	LYS
1	D	99	LEU
1	D	149	MET
1	D	161	PHE
1	D	171	MET
1	D	178	PRO
1	D	202	VAL
1	D	280	LEU
1	D	282	GLU
1	C	5	LEU
1	C	26	LEU
1	C	27	LEU
1	C	31	ARG
1	C	77	ILE
1	C	79	ARG
1	C	80	LYS
1	C	103	ILE
1	C	134	ARG
1	C	165	ASN
1	C	171	MET
1	C	175	LYS
1	C	190	LEU
1	C	197	GLU
1	C	236	ILE
1	C	237	LYS
1	C	269	LYS
1	C	283	ASP
1	A	9	LYS
1	A	27	LEU
1	A	31	ARG
1	A	53	ASP
1	A	77	ILE

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Mol	Chain	Res	Type
1	A	109	ASP
1	A	171	MET
1	A	175	LYS
1	A	190	LEU
1	A	213	ILE
1	A	298	LEU
1	A	302	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	60	ASN
1	B	165	ASN
1	B	172	HIS
1	B	174	GLN
1	D	45	HIS
1	D	60	ASN
1	D	172	HIS
1	D	174	GLN
1	D	208	ASN
1	D	251	GLN
1	D	293	GLN
1	C	45	HIS
1	C	60	ASN
1	C	165	ASN
1	C	251	GLN
1	A	165	ASN
1	A	251	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	3
1	C	2
1	A	1
1	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	219:TYR	C	220:SER	N	1.20
1	A	123:THR	C	124:TYR	N	1.20
1	B	217:ARG	C	218:GLY	N	1.19
1	B	52:VAL	C	53:ASP	N	1.18
1	B	240:SER	C	241:LYS	N	1.18
1	D	135:GLU	C	136:ARG	N	1.17
1	C	253:GLU	C	254:TYR	N	1.16

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	298/308 (96%)	-0.57	3 (1%) 82 82	21, 35, 58, 74	0
1	B	301/308 (97%)	-0.44	7 (2%) 60 59	18, 34, 53, 79	0
1	C	299/308 (97%)	-0.49	5 (1%) 70 70	23, 39, 62, 82	0
1	D	308/308 (100%)	-0.61	9 (2%) 51 48	20, 34, 59, 71	0
All	All	1206/1232 (97%)	-0.53	24 (1%) 65 63	18, 35, 59, 82	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	87	GLN	5.7
1	D	85	ARG	4.2
1	D	84	THR	4.1
1	D	87	GLN	3.2
1	B	305	GLN	3.2
1	D	86	GLU	3.0
1	D	83	MET	3.0
1	A	77	ILE	2.9
1	C	79	ARG	2.5
1	B	76	GLY	2.5
1	D	76	GLY	2.5
1	D	258	ASP	2.4
1	C	78	GLY	2.4
1	D	79	ARG	2.3
1	D	77	ILE	2.2
1	B	115	THR	2.2
1	B	78	GLY	2.2
1	C	77	ILE	2.2
1	C	115	THR	2.2
1	A	305	GLN	2.1
1	C	76	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	78	GLY	2.1
1	B	86	GLU	2.0
1	B	77	ILE	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 monosaccharides 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.