



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 22, 2021 – 02:19 PM EDT

PDB ID : 2O6Y
Title : Tyrosine ammonia-lyase from Rhodobacter sphaeroides
Authors : Louie, G.V.; Bowman, M.E.; Moffitt, M.C.; Baiga, T.J.; Moore, B.S.; Noel, J.P.
Deposited on : 2006-12-09
Resolution : 1.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.22
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) : 2.22

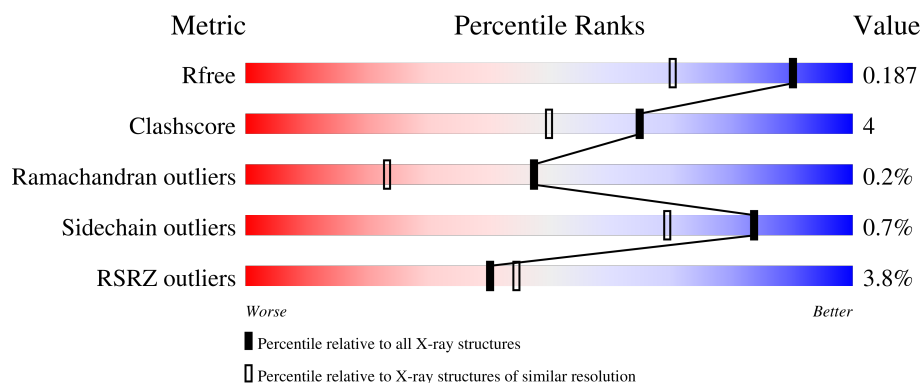
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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	521	<div> <div>5%</div> <div> <div></div> <div>91%</div> <div>7%</div> </div> </div>
1	B	521	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>11%</div> </div> </div>
1	C	521	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>9%</div> </div> </div>
1	D	521	<div> <div>5%</div> <div> <div></div> <div>88%</div> <div>10%</div> </div> </div>
1	E	521	<div> <div>4%</div> <div> <div></div> <div>91%</div> <div>7%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	521	<div><div></div><div>2%</div><div>91%</div><div>7% ..</div></div>
1	G	521	<div><div></div><div>2%</div><div>90%</div><div>8% .</div></div>
1	H	521	<div><div></div><div>6%</div><div>88%</div><div>10% .</div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 34467 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 Putative histidine ammonia-lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	515	Total	C	N	O	S	0	0	0
			3813	2363	731	705	14			
1	B	515	Total	C	N	O	S	0	0	0
			3813	2363	731	705	14			
1	C	514	Total	C	N	O	S	0	0	0
			3806	2358	730	704	14			
1	D	514	Total	C	N	O	S	0	0	0
			3806	2358	730	704	14			
1	E	514	Total	C	N	O	S	0	0	0
			3806	2358	730	704	14			
1	F	514	Total	C	N	O	S	0	0	0
			3806	2358	730	704	14			
1	G	514	Total	C	N	O	S	0	0	0
			3806	2358	730	704	14			
1	H	514	Total	C	N	O	S	0	0	0
			3806	2358	730	704	14			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
A	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
A	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
B	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
B	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
B	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
C	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
C	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
C	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
D	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
D	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
D	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
E	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0

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Chain	Residue	Modelled	Actual	Comment	Reference
E	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
E	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
F	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
F	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
F	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
G	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
G	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
G	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
H	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
H	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
H	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0

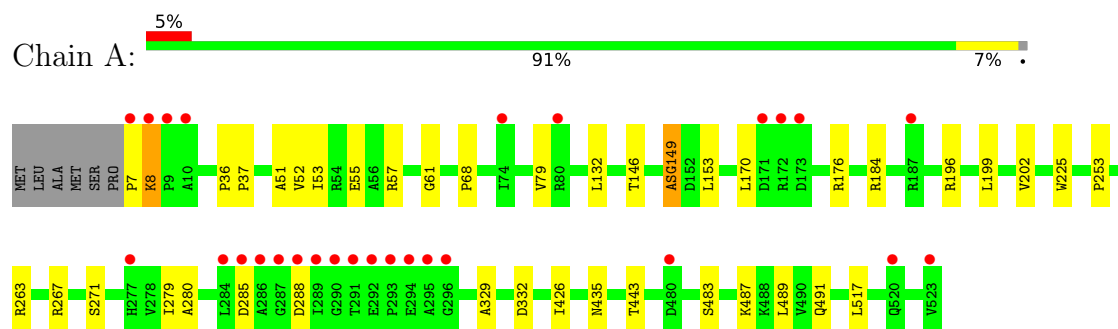
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	458	Total O 458 458	0	0
2	B	487	Total O 487 487	0	0
2	C	496	Total O 496 496	0	0
2	D	494	Total O 494 494	0	0
2	E	498	Total O 498 498	0	0
2	F	586	Total O 586 586	0	0
2	G	523	Total O 523 523	0	0
2	H	463	Total O 463 463	0	0

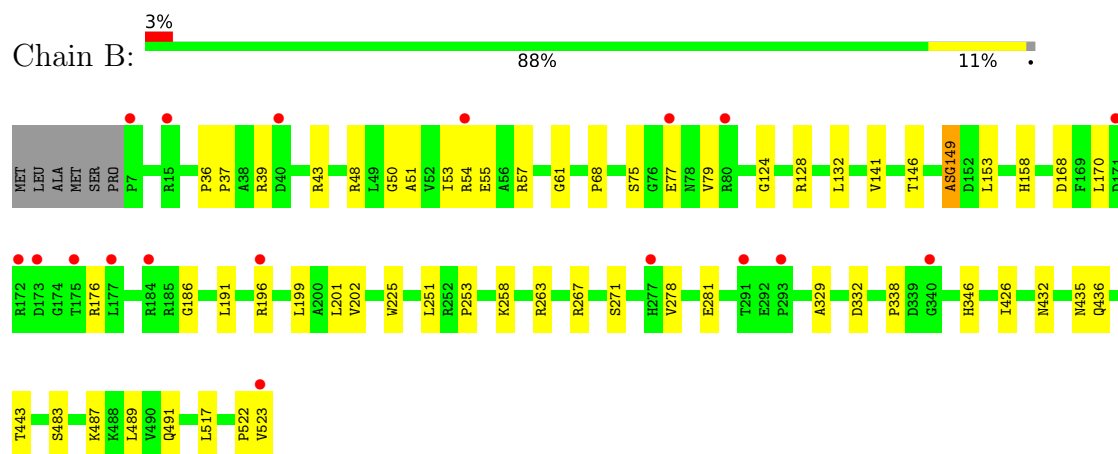
3 Residue-property plots [i](#)

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.

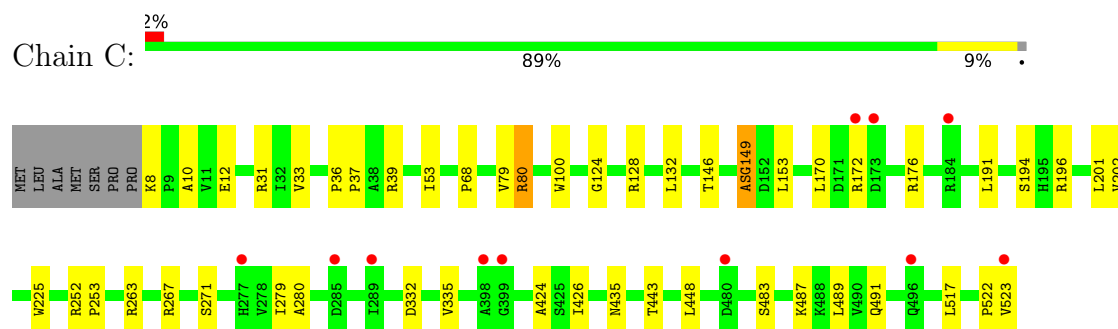
- Molecule 1: Putative histidine ammonia-lyase




- Molecule 1: Putative histidine ammonia-lyase

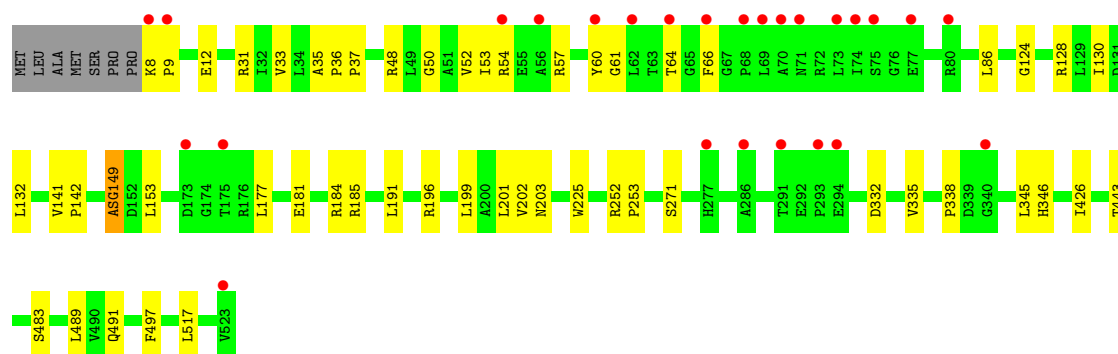


- Molecule 1: Putative histidine ammonia-lyase

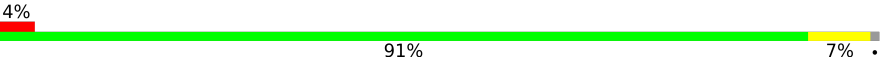


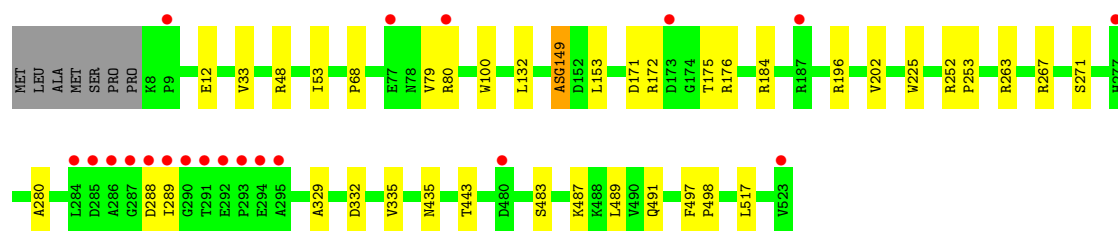
- Molecule 1: Putative histidine ammonia-lyase

Chain D: 

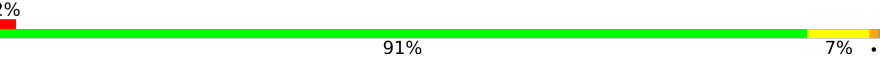


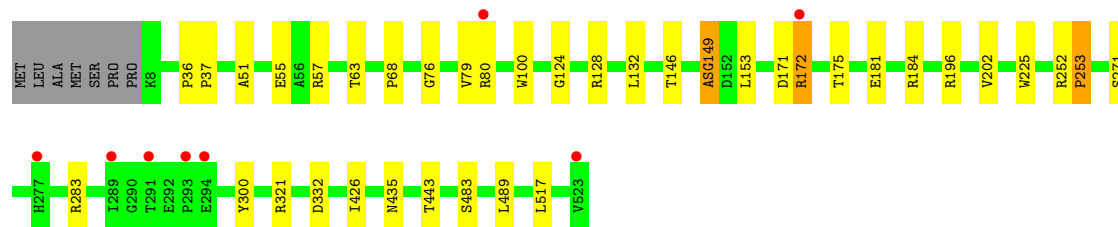
• Molecule 1: Putative histidine ammonia-lyase

Chain E: 




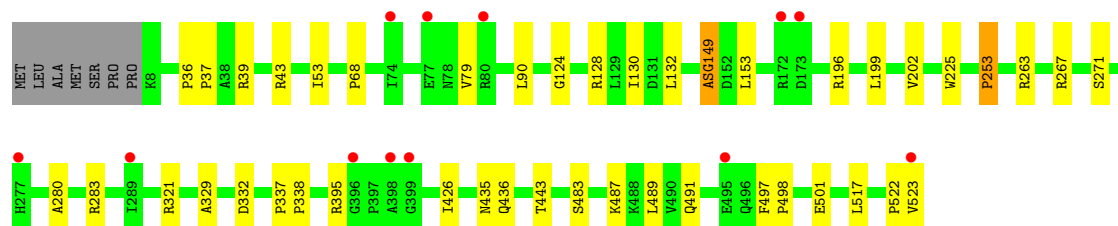
• Molecule 1: Putative histidine ammonia-lyase

Chain F: 



• Molecule 1: Putative histidine ammonia-lyase

Chain G: 



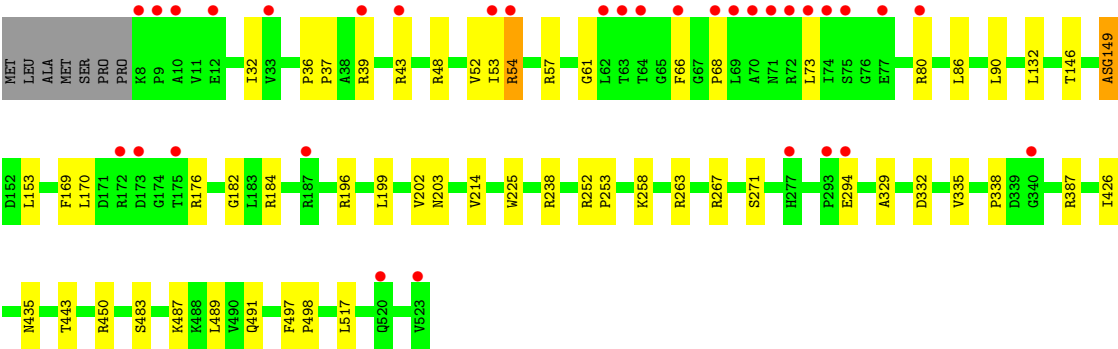
• Molecule 1: Putative histidine ammonia-lyase

Chain H:

6%

88%

10%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.92Å 155.63Å 164.35Å 90.00° 94.13° 90.00°	Depositor
Resolution (Å)	500.00 – 1.50 25.78 – 1.50	Depositor EDS
% Data completeness (in resolution range)	92.0 (500.00-1.50) 93.2 (25.78-1.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 1.50Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.174 , 0.189 0.173 , 0.187	Depositor DCC
R_{free} test set	32530 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	13.0	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	34467	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.13% 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: MDO

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	0.25	0/3862	0.54	0/5253
1	B	0.25	0/3862	0.54	0/5253
1	C	0.26	0/3854	0.55	0/5242
1	D	0.25	0/3854	0.55	0/5242
1	E	0.25	0/3854	0.55	0/5242
1	F	0.26	0/3854	0.57	1/5242 (0.0%)
1	G	0.26	0/3854	0.57	1/5242 (0.0%)
1	H	0.24	0/3854	0.54	0/5242
All	All	0.25	0/30848	0.55	2/41958 (0.0%)

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	B	0	2
1	C	0	2
1	D	0	2
1	E	0	2
1	F	0	2
1	G	0	2
1	H	0	2
All	All	0	16

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	253	PRO	N-CA-C	5.20	125.62	112.10
1	G	253	PRO	N-CA-C	5.02	125.15	112.10

There are no chirality outliers.

All (16) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	149	MDO	Peptide,Mainchain
1	B	149	MDO	Peptide,Mainchain
1	C	149	MDO	Peptide,Mainchain
1	D	149	MDO	Peptide,Mainchain
1	E	149	MDO	Peptide,Mainchain
1	F	149	MDO	Peptide,Mainchain
1	G	149	MDO	Peptide,Mainchain
1	H	149	MDO	Peptide,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	3813	0	3864	25	0
1	B	3813	0	3864	39	0
1	C	3806	0	3856	35	0
1	D	3806	0	3856	34	0
1	E	3806	0	3856	29	0
1	F	3806	0	3856	27	0
1	G	3806	0	3856	28	0
1	H	3806	0	3856	48	0
2	A	458	0	0	3	0
2	B	487	0	0	5	0
2	C	496	0	0	4	0
2	D	494	0	0	4	0
2	E	498	0	0	7	0
2	F	586	0	0	6	0
2	G	523	0	0	4	0
2	H	463	0	0	5	0
All	All	34467	0	30864	250	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 4.

All (250) 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:D:153:LEU:HD22	1:D:202:VAL:HG12	1.63	0.80
1:B:53:ILE:HG21	1:B:196:ARG:HH11	1.44	0.80
1:H:53:ILE:HG21	1:H:196:ARG:HH11	1.47	0.79
1:H:153:LEU:HD22	1:H:202:VAL:HG12	1.65	0.78
1:E:153:LEU:HD22	1:E:202:VAL:HG12	1.70	0.73
1:E:487:LYS:O	1:E:491:GLN:HG3	1.88	0.73
1:B:153:LEU:HD22	1:B:202:VAL:HG12	1.69	0.73
1:B:487:LYS:O	1:B:491:GLN:HG3	1.89	0.73
1:A:153:LEU:HD22	1:A:202:VAL:HG12	1.70	0.73
1:D:61:GLY:HA3	1:D:199:LEU:HD11	1.70	0.71
1:F:153:LEU:HD22	1:F:202:VAL:HG12	1.72	0.70
1:C:487:LYS:O	1:C:491:GLN:HG3	1.92	0.69
1:E:172:ARG:NH1	2:E:4898:HOH:O	2.24	0.69
1:H:238:ARG:HH11	1:H:387:ARG:NH2	1.91	0.69
1:F:79:VAL:HG11	1:F:196:ARG:HD3	1.76	0.68
1:C:153:LEU:HD22	1:C:202:VAL:HG12	1.76	0.67
1:D:53:ILE:HD13	1:D:196:ARG:HD2	1.78	0.64
1:H:238:ARG:NH1	1:H:387:ARG:NH2	2.47	0.63
1:H:53:ILE:HG21	1:H:196:ARG:NH1	2.14	0.63
1:D:50:GLY:O	1:D:54:ARG:HG3	1.99	0.62
1:H:487:LYS:O	1:H:491:GLN:HG3	2.00	0.62
1:H:214:VAL:HB	1:H:450:ARG:NH2	2.16	0.61
1:F:68:PRO:HG3	1:F:435:ASN:HB2	1.84	0.60
1:G:283:ARG:HD3	2:G:2777:HOH:O	2.00	0.60
1:B:53:ILE:CG2	1:B:196:ARG:HH11	2.14	0.60
1:G:53:ILE:HD13	1:G:196:ARG:HD2	1.83	0.60
1:G:153:LEU:HD22	1:G:202:VAL:HG12	1.83	0.60
1:G:487:LYS:O	1:G:491:GLN:HG3	2.02	0.60
1:B:51:ALA:O	1:B:55:GLU:HG3	2.02	0.59
1:B:79:VAL:HG21	2:B:4884:HOH:O	2.03	0.58
1:B:68:PRO:HG3	1:B:435:ASN:HB2	1.85	0.57
1:F:283:ARG:HD3	2:F:3716:HOH:O	2.04	0.57
1:F:171:ASP:OD2	1:F:175:THR:HB	2.05	0.57
1:H:39:ARG:HG3	1:H:39:ARG:HH11	1.71	0.56
1:E:184:ARG:HD2	2:E:3810:HOH:O	2.06	0.56
1:A:271:SER:HB2	1:A:483:SER:HB3	1.88	0.55
1:H:184:ARG:HG2	1:H:184:ARG:HH11	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:LEU:HD23	1:A:176:ARG:HG2	1.88	0.54
1:D:52:VAL:HG13	1:D:57:ARG:HB2	1.88	0.54
1:B:432:ASN:HB2	1:B:436:GLN:HE21	1.72	0.54
1:E:271:SER:HB2	1:E:483:SER:HB3	1.90	0.54
1:C:124:GLY:O	1:C:128:ARG:HG2	2.08	0.54
1:B:61:GLY:HA3	1:B:199:LEU:HD11	1.89	0.53
1:C:176:ARG:HD2	2:C:3285:HOH:O	2.08	0.53
1:H:54:ARG:HG2	1:H:54:ARG:HH11	1.73	0.53
1:H:271:SER:HB2	1:H:483:SER:HB3	1.90	0.53
1:D:12:GLU:HA	1:D:33:VAL:HG13	1.90	0.53
1:D:36:PRO:HB2	1:D:37:PRO:HD3	1.89	0.53
1:F:124:GLY:O	1:F:128:ARG:HG2	2.09	0.53
1:G:271:SER:HB2	1:G:483:SER:HB3	1.90	0.53
1:H:61:GLY:HA3	1:H:199:LEU:HD21	1.90	0.53
1:B:53:ILE:HG21	1:B:196:ARG:NH1	2.21	0.53
1:E:225:TRP:CE2	1:E:517:LEU:HD22	2.44	0.53
1:A:279:ILE:HD11	1:D:346:HIS:CE1	2.44	0.52
1:E:79:VAL:HG21	2:E:2045:HOH:O	2.09	0.52
2:A:3436:HOH:O	1:D:426:ILE:HD13	2.09	0.52
1:G:523:VAL:OXT	1:G:523:VAL:HG12	2.08	0.52
1:A:487:LYS:O	1:A:491:GLN:HG3	2.10	0.52
1:C:79:VAL:HG11	1:C:196:ARG:HD3	1.92	0.52
1:C:12:GLU:HA	1:C:33:VAL:HG13	1.92	0.52
1:H:225:TRP:CE2	1:H:517:LEU:HD22	2.45	0.51
1:A:225:TRP:CE2	1:A:517:LEU:HD22	2.46	0.51
2:B:3548:HOH:O	1:C:426:ILE:HD13	2.10	0.51
1:F:426:ILE:HD13	2:G:3571:HOH:O	2.10	0.51
1:H:68:PRO:HG3	1:H:435:ASN:HB2	1.92	0.51
1:B:53:ILE:HD13	1:B:196:ARG:HD2	1.93	0.51
1:A:426:ILE:HD13	2:D:3398:HOH:O	2.10	0.51
1:B:191:LEU:HD21	1:B:201:LEU:CD1	2.40	0.51
1:A:53:ILE:HD13	1:A:196:ARG:HD2	1.92	0.51
1:E:196:ARG:HG3	1:E:196:ARG:HH11	1.75	0.50
1:B:168:ASP:HB2	1:B:176:ARG:HH21	1.77	0.50
1:B:124:GLY:O	1:B:128:ARG:HG2	2.11	0.50
1:C:80:ARG:HE	1:C:194:SER:HA	1.77	0.50
1:A:176:ARG:HD2	2:A:3332:HOH:O	2.11	0.50
1:B:258:LYS:HD2	2:B:3514:HOH:O	2.12	0.50
1:A:170:LEU:CD2	1:A:176:ARG:HG2	2.42	0.50
1:B:251:LEU:HD13	1:C:335:VAL:HG21	1.93	0.50
1:D:66:PHE:CE2	1:D:86:LEU:HB2	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:184:ARG:HG2	1:H:184:ARG:NH1	2.28	0.49
1:C:271:SER:HB2	1:C:483:SER:HB3	1.94	0.49
1:A:61:GLY:HA3	1:A:199:LEU:HD21	1.95	0.49
1:F:225:TRP:CE2	1:F:517:LEU:HD22	2.47	0.49
1:B:36:PRO:HB2	1:B:37:PRO:HD3	1.93	0.49
1:B:39:ARG:HD2	1:B:43:ARG:NH2	2.28	0.49
1:H:66:PHE:CZ	1:H:86:LEU:HD22	2.48	0.49
1:E:263:ARG:O	1:E:267:ARG:HG2	2.13	0.48
1:E:132:LEU:C	1:E:132:LEU:HD13	2.32	0.48
1:A:132:LEU:C	1:A:132:LEU:HD13	2.33	0.48
1:A:263:ARG:O	1:A:267:ARG:HG2	2.13	0.48
1:A:280:ALA:O	1:D:57:ARG:HD2	2.13	0.48
1:C:225:TRP:CE2	1:C:517:LEU:HD22	2.48	0.48
1:F:76:GLY:O	1:F:79:VAL:HG23	2.13	0.48
1:A:68:PRO:HG3	1:A:435:ASN:HB2	1.95	0.48
1:D:184:ARG:HG2	1:D:184:ARG:HH11	1.78	0.48
1:D:225:TRP:CE2	1:D:517:LEU:HD22	2.49	0.48
1:E:12:GLU:HA	1:E:33:VAL:HG13	1.95	0.48
1:H:90:LEU:HD13	1:H:153:LEU:HB3	1.96	0.48
1:C:53:ILE:HD13	1:C:196:ARG:HD2	1.94	0.48
1:A:51:ALA:O	1:A:55:GLU:HG3	2.13	0.48
1:C:39:ARG:NH1	2:C:4878:HOH:O	2.46	0.48
1:B:141:VAL:HG13	1:B:158:HIS:HB2	1.96	0.48
1:C:68:PRO:HG3	1:C:435:ASN:HB2	1.96	0.48
1:G:39:ARG:HD2	1:G:43:ARG:NH2	2.29	0.48
1:E:176:ARG:HD2	2:E:3308:HOH:O	2.13	0.48
1:F:79:VAL:CG1	1:F:196:ARG:HD3	2.44	0.47
1:G:124:GLY:O	1:G:128:ARG:HG2	2.14	0.47
1:B:278:VAL:HB	1:B:281:GLU:HG3	1.96	0.47
1:C:80:ARG:NE	1:C:194:SER:HA	2.29	0.47
1:G:130:ILE:HD12	2:G:4129:HOH:O	2.13	0.47
1:B:57:ARG:HD2	1:C:280:ALA:O	2.15	0.47
1:H:294:GLU:OE1	1:H:387:ARG:NH2	2.48	0.47
1:B:75:SER:OG	1:B:77:GLU:HG2	2.14	0.47
1:B:523:VAL:OXT	1:B:523:VAL:HG12	2.14	0.47
1:C:36:PRO:HB2	1:C:37:PRO:HD3	1.96	0.47
1:H:52:VAL:HG13	1:H:57:ARG:HB2	1.97	0.47
1:H:450:ARG:HG2	1:H:450:ARG:HH11	1.79	0.47
1:B:225:TRP:CE2	1:B:517:LEU:HD22	2.50	0.47
1:D:497:PHE:HB3	2:D:4227:HOH:O	2.13	0.47
1:A:79:VAL:HG11	1:A:196:ARG:HD3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:ARG:HH21	1:C:194:SER:CB	2.28	0.46
1:G:225:TRP:CE2	1:G:517:LEU:HD22	2.50	0.46
1:B:329:ALA:HB1	1:C:252:ARG:HA	1.97	0.46
1:F:36:PRO:HB2	1:F:37:PRO:HD3	1.98	0.46
1:F:80:ARG:HG2	2:F:2717:HOH:O	2.14	0.46
2:F:3660:HOH:O	1:G:426:ILE:HD13	2.14	0.46
1:H:36:PRO:N	1:H:37:PRO:HD2	2.29	0.46
1:A:36:PRO:N	1:A:37:PRO:HD2	2.30	0.46
1:E:48:ARG:HH11	1:E:48:ARG:HG2	1.80	0.46
1:B:50:GLY:O	1:B:54:ARG:HG3	2.16	0.46
1:G:132:LEU:HD13	1:G:132:LEU:C	2.36	0.46
1:H:170:LEU:HD23	1:H:176:ARG:HG2	1.98	0.46
1:E:80:ARG:HG2	2:E:3706:HOH:O	2.16	0.46
1:D:132:LEU:HD13	1:D:132:LEU:C	2.35	0.46
1:G:395:ARG:CZ	1:G:501:GLU:HG2	2.46	0.46
1:B:271:SER:HB2	1:B:483:SER:HB3	1.98	0.45
1:D:12:GLU:HG2	1:D:35:ALA:HB2	1.98	0.45
1:F:271:SER:HB2	1:F:483:SER:HB3	1.98	0.45
1:H:132:LEU:C	1:H:132:LEU:HD13	2.37	0.45
1:A:7:PRO:O	1:A:8:LYS:C	2.55	0.45
1:B:132:LEU:HD13	1:B:132:LEU:C	2.37	0.45
1:C:170:LEU:HD23	1:C:176:ARG:HG2	1.98	0.45
1:C:225:TRP:CE2	1:C:522:PRO:HG3	2.52	0.45
1:A:329:ALA:HB1	1:D:252:ARG:HA	1.99	0.45
1:B:170:LEU:CD2	1:B:176:ARG:HG2	2.46	0.45
1:E:100:TRP:CZ2	1:E:172:ARG:HG2	2.52	0.45
1:C:132:LEU:C	1:C:132:LEU:HD13	2.37	0.45
1:E:171:ASP:OD2	1:E:175:THR:HB	2.15	0.45
1:F:252:ARG:HA	1:G:329:ALA:HB1	1.97	0.45
1:A:52:VAL:HG13	1:A:57:ARG:HB2	1.99	0.45
1:E:288:ASP:O	1:H:73:LEU:HD12	2.16	0.45
1:H:170:LEU:CD2	1:H:176:ARG:HG2	2.47	0.45
1:H:176:ARG:HD2	2:H:3379:HOH:O	2.16	0.44
1:B:346:HIS:CE1	1:C:279:ILE:HD11	2.52	0.44
1:B:426:ILE:HD13	2:C:3493:HOH:O	2.16	0.44
1:H:39:ARG:O	1:H:43:ARG:HG3	2.16	0.44
1:D:48:ARG:NH2	1:D:338:PRO:O	2.50	0.44
1:C:523:VAL:OXT	1:C:523:VAL:HG12	2.18	0.44
1:G:36:PRO:N	1:G:37:PRO:HD2	2.33	0.44
1:H:238:ARG:NH1	1:H:387:ARG:HH22	2.14	0.44
1:D:489:LEU:C	1:D:489:LEU:HD23	2.37	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:146:THR:HG22	2:H:2163:HOH:O	2.18	0.44
1:H:39:ARG:HD3	1:H:43:ARG:HH21	1.83	0.44
1:H:258:LYS:HD2	2:H:3845:HOH:O	2.16	0.44
1:H:335:VAL:HG11	2:H:1982:HOH:O	2.17	0.44
1:A:489:LEU:C	1:A:489:LEU:HD23	2.38	0.44
1:B:225:TRP:CE2	1:B:522:PRO:HD3	2.53	0.44
1:G:79:VAL:HG11	1:G:196:ARG:HD3	1.99	0.44
1:C:12:GLU:HA	1:C:33:VAL:CG1	2.47	0.44
1:E:280:ALA:O	1:H:57:ARG:HD2	2.17	0.44
1:E:489:LEU:HD23	1:E:489:LEU:C	2.37	0.44
1:E:196:ARG:HG3	1:E:196:ARG:NH1	2.32	0.43
1:G:68:PRO:HG3	1:G:435:ASN:HB2	2.00	0.43
1:H:214:VAL:HB	1:H:450:ARG:HH21	1.82	0.43
1:D:177:LEU:HD21	1:D:185:ARG:NH1	2.33	0.43
1:D:191:LEU:HD21	1:D:201:LEU:CD1	2.49	0.43
1:H:53:ILE:HD13	1:H:196:ARG:HD2	1.99	0.43
1:B:489:LEU:HD23	1:B:489:LEU:C	2.38	0.43
1:C:263:ARG:O	1:C:267:ARG:HG2	2.18	0.43
1:D:271:SER:HB2	1:D:483:SER:HB3	2.00	0.43
1:F:132:LEU:C	1:F:132:LEU:HD13	2.38	0.43
1:F:146:THR:HG22	2:F:2088:HOH:O	2.19	0.43
1:F:171:ASP:C	1:F:171:ASP:OD1	2.57	0.43
1:C:489:LEU:C	1:C:489:LEU:HD23	2.38	0.43
1:H:199:LEU:HD12	1:H:199:LEU:N	2.33	0.43
1:B:48:ARG:NH2	1:B:338:PRO:O	2.51	0.43
1:D:8:LYS:CG	1:D:31:ARG:HG3	2.48	0.43
1:D:66:PHE:CZ	1:D:86:LEU:HD22	2.53	0.43
1:H:199:LEU:O	1:H:203:ASN:HB2	2.19	0.43
1:H:263:ARG:O	1:H:267:ARG:HG2	2.18	0.43
1:D:335:VAL:HG23	1:D:345:LEU:HB2	2.01	0.43
2:E:3348:HOH:O	1:H:426:ILE:HD13	2.18	0.43
1:F:63:THR:HG22	2:F:4294:HOH:O	2.17	0.43
1:G:225:TRP:CE2	1:G:522:PRO:HD3	2.54	0.43
1:B:186:GLY:HA2	2:B:3021:HOH:O	2.19	0.43
1:F:100:TRP:CH2	1:F:172:ARG:HD3	2.54	0.43
1:H:66:PHE:CE2	1:H:86:LEU:HB2	2.54	0.43
1:D:124:GLY:O	1:D:128:ARG:HG2	2.19	0.43
1:E:289:ILE:HD11	2:H:4360:HOH:O	2.18	0.43
1:F:489:LEU:C	1:F:489:LEU:HD23	2.39	0.43
1:F:181:GLU:OE1	1:F:184:ARG:NH1	2.52	0.43
1:F:321:ARG:HH21	1:G:321:ARG:HH21	1.65	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:169:PHE:CD2	1:H:182:GLY:HA3	2.54	0.43
1:D:199:LEU:O	1:D:203:ASN:HB2	2.19	0.42
1:E:48:ARG:HG2	1:E:48:ARG:NH1	2.34	0.42
1:E:335:VAL:HG11	2:E:1376:HOH:O	2.19	0.42
1:H:80:ARG:HH11	1:H:80:ARG:HG3	1.84	0.42
1:D:184:ARG:HG2	1:D:184:ARG:NH1	2.33	0.42
1:E:252:ARG:HA	1:H:329:ALA:HB1	2.01	0.42
1:F:79:VAL:HG21	2:F:2178:HOH:O	2.19	0.42
1:C:146:THR:HG22	2:C:1832:HOH:O	2.19	0.42
1:B:170:LEU:HD23	1:B:176:ARG:HG2	2.01	0.42
1:D:130:ILE:HD12	2:D:4714:HOH:O	2.19	0.42
1:G:489:LEU:HD23	1:G:489:LEU:C	2.39	0.42
1:C:225:TRP:CD2	1:C:522:PRO:HG3	2.54	0.42
1:C:100:TRP:CZ2	1:C:172:ARG:HG2	2.55	0.42
1:D:491:GLN:NE2	2:D:2981:HOH:O	2.52	0.42
1:F:68:PRO:HG3	1:F:435:ASN:CB	2.49	0.42
1:B:263:ARG:O	1:B:267:ARG:HG2	2.19	0.41
1:G:497:PHE:HA	1:G:498:PRO:HD2	1.94	0.41
1:C:8:LYS:HE3	1:C:8:LYS:HB2	1.92	0.41
1:H:48:ARG:NH2	1:H:338:PRO:O	2.53	0.41
1:H:497:PHE:HA	1:H:498:PRO:HD3	1.91	0.41
1:F:57:ARG:HD2	1:G:280:ALA:O	2.20	0.41
1:H:48:ARG:O	1:H:52:VAL:HG23	2.19	0.41
1:A:146:THR:HG22	2:A:2274:HOH:O	2.20	0.41
1:H:489:LEU:HD23	1:H:489:LEU:C	2.40	0.41
1:E:68:PRO:HG3	1:E:435:ASN:HB2	2.01	0.41
1:G:199:LEU:HD23	2:G:4049:HOH:O	2.19	0.41
1:E:53:ILE:HD13	1:E:196:ARG:HD2	2.02	0.41
1:F:51:ALA:O	1:F:55:GLU:HG3	2.20	0.41
1:G:395:ARG:NH1	1:G:501:GLU:HG2	2.35	0.41
1:B:146:THR:HG22	2:B:2492:HOH:O	2.21	0.41
1:F:300:TYR:CZ	1:G:436:GLN:HG2	2.56	0.41
1:G:90:LEU:HD13	1:G:153:LEU:HB3	2.01	0.41
1:C:191:LEU:HD21	1:C:201:LEU:CD1	2.51	0.41
1:A:285:ASP:OD1	1:A:288:ASP:OD2	2.39	0.41
1:B:53:ILE:CG2	1:B:196:ARG:NH1	2.82	0.41
1:C:424:ALA:HB2	1:C:448:LEU:HD12	2.03	0.41
1:D:60:TYR:HA	1:D:64:THR:OG1	2.21	0.41
1:D:181:GLU:OE1	1:D:181:GLU:HA	2.21	0.41
1:G:337:PRO:HA	1:G:338:PRO:HD3	1.96	0.41
1:A:184:ARG:HG2	1:A:184:ARG:HH11	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:12:GLU:HA	1:E:33:VAL:CG1	2.51	0.40
1:D:141:VAL:HA	1:D:142:PRO:HD3	1.99	0.40
1:G:263:ARG:O	1:G:267:ARG:HG2	2.21	0.40
1:C:10:ALA:HB2	1:C:31:ARG:HH21	1.87	0.40
1:E:329:ALA:HB1	1:H:252:ARG:HA	2.03	0.40
1:E:497:PHE:HA	1:E:498:PRO:HD3	1.91	0.40
1:C:39:ARG:NH1	1:C:39:ARG:HG2	2.36	0.40
1:D:8:LYS:HB2	1:D:9:PRO:HD2	2.04	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	512/521 (98%)	504 (98%)	6 (1%)	2 (0%)	34	13
1	B	512/521 (98%)	507 (99%)	4 (1%)	1 (0%)	47	23
1	C	511/521 (98%)	506 (99%)	4 (1%)	1 (0%)	47	23
1	D	511/521 (98%)	504 (99%)	6 (1%)	1 (0%)	47	23
1	E	511/521 (98%)	504 (99%)	6 (1%)	1 (0%)	47	23
1	F	511/521 (98%)	504 (99%)	6 (1%)	1 (0%)	47	23
1	G	511/521 (98%)	506 (99%)	4 (1%)	1 (0%)	47	23
1	H	511/521 (98%)	504 (99%)	6 (1%)	1 (0%)	47	23
All	All	4090/4168 (98%)	4039 (99%)	42 (1%)	9 (0%)	47	23

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	253	PRO

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Mol	Chain	Res	Type
1	E	253	PRO
1	B	253	PRO
1	C	253	PRO
1	D	253	PRO
1	F	253	PRO
1	G	253	PRO
1	H	253	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	384/389 (99%)	382 (100%)	2 (0%)	88	78
1	B	384/389 (99%)	382 (100%)	2 (0%)	88	78
1	C	383/389 (98%)	380 (99%)	3 (1%)	81	66
1	D	383/389 (98%)	381 (100%)	2 (0%)	88	78
1	E	383/389 (98%)	381 (100%)	2 (0%)	88	78
1	F	383/389 (98%)	380 (99%)	3 (1%)	81	66
1	G	383/389 (98%)	381 (100%)	2 (0%)	88	78
1	H	383/389 (98%)	379 (99%)	4 (1%)	76	57
All	All	3066/3112 (98%)	3046 (99%)	20 (1%)	84	69

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	332	ASP
1	A	443	THR
1	B	332	ASP
1	B	443	THR
1	C	80	ARG
1	C	332	ASP
1	C	443	THR
1	D	332	ASP

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Mol	Chain	Res	Type
1	D	443	THR
1	E	332	ASP
1	E	443	THR
1	F	172	ARG
1	F	332	ASP
1	F	443	THR
1	G	332	ASP
1	G	443	THR
1	H	32	ILE
1	H	54	ARG
1	H	332	ASP
1	H	443	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	394	HIS
1	A	491	GLN
1	B	394	HIS
1	B	436	GLN
1	C	394	HIS
1	D	189	GLN
1	D	394	HIS
1	D	491	GLN
1	D	496	GLN
1	E	189	GLN
1	E	394	HIS
1	E	520	GLN
1	F	394	HIS
1	F	491	GLN
1	G	394	HIS
1	H	78	ASN
1	H	394	HIS
1	H	508	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	MDO	E	149	1	12,13,14	2.20	4 (33%)	15,18,20	4.13	6 (40%)
1	MDO	G	149	1	12,13,14	2.19	4 (33%)	15,18,20	4.16	6 (40%)
1	MDO	C	149	1	12,13,14	2.27	4 (33%)	15,18,20	4.07	6 (40%)
1	MDO	B	149	1	12,13,14	2.24	4 (33%)	15,18,20	3.93	6 (40%)
1	MDO	H	149	1	12,13,14	2.26	4 (33%)	15,18,20	4.03	6 (40%)
1	MDO	F	149	1	12,13,14	2.20	4 (33%)	15,18,20	4.04	6 (40%)
1	MDO	A	149	1	12,13,14	2.23	4 (33%)	15,18,20	4.03	6 (40%)
1	MDO	D	149	1	12,13,14	2.22	4 (33%)	15,18,20	4.04	6 (40%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MDO	E	149	1	-	0/4/23/24	0/1/1/1
1	MDO	G	149	1	-	2/4/23/24	0/1/1/1
1	MDO	C	149	1	-	0/4/23/24	0/1/1/1
1	MDO	B	149	1	-	0/4/23/24	0/1/1/1
1	MDO	H	149	1	-	0/4/23/24	0/1/1/1
1	MDO	F	149	1	-	1/4/23/24	0/1/1/1
1	MDO	A	149	1	-	0/4/23/24	0/1/1/1
1	MDO	D	149	1	-	0/4/23/24	0/1/1/1

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	149	MDO	O2-C2	5.00	1.33	1.23
1	B	149	MDO	O2-C2	4.95	1.33	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	149	MDO	O2-C2	4.90	1.33	1.23
1	D	149	MDO	O2-C2	4.78	1.33	1.23
1	A	149	MDO	O2-C2	4.78	1.33	1.23
1	F	149	MDO	O2-C2	4.68	1.33	1.23
1	E	149	MDO	O2-C2	4.68	1.33	1.23
1	G	149	MDO	O2-C2	4.56	1.32	1.23
1	B	149	MDO	C2-N3	-3.67	1.31	1.39
1	C	149	MDO	C2-N3	-3.66	1.31	1.39
1	G	149	MDO	C2-N3	-3.65	1.31	1.39
1	H	149	MDO	C2-N3	-3.64	1.31	1.39
1	E	149	MDO	C2-N3	-3.61	1.31	1.39
1	A	149	MDO	C2-N3	-3.58	1.31	1.39
1	F	149	MDO	C2-N3	-3.49	1.31	1.39
1	D	149	MDO	C2-N3	-3.48	1.31	1.39
1	A	149	MDO	CA2-N2	-2.71	1.34	1.39
1	C	149	MDO	CA2-N2	-2.71	1.34	1.39
1	D	149	MDO	CA2-N2	-2.65	1.34	1.39
1	E	149	MDO	CA2-N2	-2.61	1.34	1.39
1	G	149	MDO	CA2-N2	-2.60	1.34	1.39
1	B	149	MDO	CA2-N2	-2.60	1.34	1.39
1	F	149	MDO	CA2-N2	-2.59	1.34	1.39
1	H	149	MDO	CA2-N2	-2.58	1.34	1.39
1	H	149	MDO	CA2-C2	-2.57	1.38	1.43
1	E	149	MDO	CA2-C2	-2.57	1.38	1.43
1	F	149	MDO	CA2-C2	-2.50	1.38	1.43
1	G	149	MDO	CA2-C2	-2.49	1.38	1.43
1	A	149	MDO	CA2-C2	-2.47	1.38	1.43
1	D	149	MDO	CA2-C2	-2.42	1.38	1.43
1	C	149	MDO	CA2-C2	-2.35	1.38	1.43
1	B	149	MDO	CA2-C2	-2.30	1.38	1.43

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	149	MDO	CA2-C2-N3	13.22	109.62	103.37
1	E	149	MDO	CA2-C2-N3	13.13	109.58	103.37
1	H	149	MDO	CA2-C2-N3	12.95	109.49	103.37
1	C	149	MDO	CA2-C2-N3	12.92	109.48	103.37
1	D	149	MDO	CA2-C2-N3	12.87	109.45	103.37
1	A	149	MDO	CA2-C2-N3	12.78	109.42	103.37
1	F	149	MDO	CA2-C2-N3	12.76	109.41	103.37
1	B	149	MDO	CA2-C2-N3	12.49	109.28	103.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	149	MDO	O2-C2-CA2	-5.19	128.04	130.96
1	E	149	MDO	O2-C2-CA2	-5.15	128.07	130.96
1	F	149	MDO	O2-C2-CA2	-5.11	128.09	130.96
1	A	149	MDO	O2-C2-CA2	-4.92	128.19	130.96
1	C	149	MDO	O2-C2-CA2	-4.82	128.25	130.96
1	D	149	MDO	O2-C2-CA2	-4.77	128.28	130.96
1	B	149	MDO	O2-C2-CA2	-4.52	128.42	130.96
1	H	149	MDO	O2-C2-CA2	-4.49	128.44	130.96
1	B	149	MDO	C2-CA2-N2	-4.04	106.10	108.93
1	E	149	MDO	C2-CA2-N2	-4.02	106.12	108.93
1	C	149	MDO	C2-CA2-N2	-3.98	106.15	108.93
1	G	149	MDO	C2-CA2-N2	-3.97	106.15	108.93
1	F	149	MDO	C2-CA2-N2	-3.94	106.17	108.93
1	D	149	MDO	C2-CA2-N2	-3.90	106.20	108.93
1	A	149	MDO	C2-CA2-N2	-3.89	106.21	108.93
1	H	149	MDO	C2-CA2-N2	-3.86	106.23	108.93
1	C	149	MDO	CB2-CA2-C2	3.65	129.45	122.76
1	B	149	MDO	CB2-CA2-C2	3.58	129.31	122.76
1	H	149	MDO	CB2-CA2-C2	3.55	129.25	122.76
1	E	149	MDO	CB2-CA2-C2	3.49	129.15	122.76
1	A	149	MDO	CB2-CA2-C2	3.48	129.13	122.76
1	D	149	MDO	CB2-CA2-C2	3.47	129.12	122.76
1	F	149	MDO	CB2-CA2-C2	3.46	129.10	122.76
1	G	149	MDO	CB2-CA2-C2	3.46	129.09	122.76
1	G	149	MDO	N3-C1-N2	-3.08	109.32	111.45
1	H	149	MDO	N3-C1-N2	-3.05	109.34	111.45
1	C	149	MDO	N3-C1-N2	-2.99	109.39	111.45
1	D	149	MDO	N3-C1-N2	-2.99	109.39	111.45
1	A	149	MDO	N3-C1-N2	-2.98	109.39	111.45
1	F	149	MDO	N3-C1-N2	-2.95	109.41	111.45
1	E	149	MDO	N3-C1-N2	-2.81	109.51	111.45
1	B	149	MDO	N3-C1-N2	-2.81	109.51	111.45
1	C	149	MDO	CA2-N2-C1	2.66	107.80	105.40
1	A	149	MDO	CA2-N2-C1	2.62	107.77	105.40
1	G	149	MDO	CA2-N2-C1	2.59	107.74	105.40
1	D	149	MDO	CA2-N2-C1	2.58	107.73	105.40
1	E	149	MDO	CA2-N2-C1	2.57	107.72	105.40
1	F	149	MDO	CA2-N2-C1	2.57	107.72	105.40
1	B	149	MDO	CA2-N2-C1	2.55	107.71	105.40
1	H	149	MDO	CA2-N2-C1	2.47	107.64	105.40

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	F	149	MDO	N2-C1-CA-CB
1	G	149	MDO	N2-C1-CA-CB
1	G	149	MDO	N3-C1-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

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	514/521 (98%)	0.26	27 (5%) 26 29	8, 16, 34, 53	0
1	B	514/521 (98%)	0.11	18 (3%) 44 48	7, 15, 31, 48	0
1	C	513/521 (98%)	-0.03	11 (2%) 63 68	7, 13, 28, 41	0
1	D	513/521 (98%)	0.13	26 (5%) 28 30	6, 14, 34, 48	0
1	E	513/521 (98%)	0.10	20 (3%) 39 44	7, 14, 31, 59	0
1	F	513/521 (98%)	-0.14	8 (1%) 72 77	6, 11, 25, 37	0
1	G	513/521 (98%)	-0.06	12 (2%) 60 65	6, 12, 26, 39	0
1	H	513/521 (98%)	0.27	33 (6%) 19 20	7, 16, 36, 52	0
All	All	4106/4168 (98%)	0.08	155 (3%) 40 44	6, 14, 31, 59	0

All (155) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	74	ILE	11.4
1	A	286	ALA	10.7
1	A	7	PRO	10.2
1	A	523	VAL	8.7
1	H	74	ILE	8.6
1	H	523	VAL	8.1
1	G	523	VAL	8.0
1	B	523	VAL	7.7
1	A	291	THR	7.7
1	B	7	PRO	7.5
1	A	296	GLY	7.4
1	E	523	VAL	7.1
1	A	289	ILE	7.0
1	A	285	ASP	7.0
1	E	293	PRO	7.0
1	E	286	ALA	7.0

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Mol	Chain	Res	Type	RSRZ
1	C	523	VAL	6.8
1	E	288	ASP	6.7
1	F	523	VAL	6.5
1	A	293	PRO	6.4
1	D	523	VAL	6.3
1	E	285	ASP	6.2
1	H	77	GLU	5.9
1	E	291	THR	5.7
1	E	289	ILE	5.7
1	A	290	GLY	5.6
1	E	294	GLU	5.4
1	A	287	GLY	5.1
1	E	290	GLY	5.1
1	A	294	GLU	5.0
1	F	293	PRO	4.9
1	H	9	PRO	4.9
1	A	288	ASP	4.6
1	D	77	GLU	4.5
1	D	293	PRO	4.4
1	B	172	ARG	4.4
1	E	287	GLY	4.4
1	D	70	ALA	4.4
1	H	66	PHE	4.2
1	H	75	SER	4.1
1	G	399	GLY	4.0
1	B	175	THR	3.9
1	C	399	GLY	3.9
1	H	277	HIS	3.8
1	B	277	HIS	3.8
1	B	293	PRO	3.6
1	H	294	GLU	3.6
1	H	187	ARG	3.5
1	E	284	LEU	3.4
1	H	8	LYS	3.4
1	A	187	ARG	3.4
1	A	10	ALA	3.4
1	F	289	ILE	3.4
1	E	277	HIS	3.4
1	A	292	GLU	3.3
1	D	64	THR	3.3
1	A	173	ASP	3.3
1	H	70	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
1	H	62	LEU	3.2
1	D	277	HIS	3.2
1	G	398	ALA	3.2
1	H	172	ARG	3.1
1	B	80	ARG	3.1
1	D	54	ARG	3.1
1	G	80	ARG	3.0
1	C	172	ARG	3.0
1	D	9	PRO	3.0
1	H	64	THR	3.0
1	E	187	ARG	3.0
1	H	173	ASP	3.0
1	H	53	ILE	2.9
1	H	69	LEU	2.9
1	H	80	ARG	2.9
1	D	80	ARG	2.9
1	F	291	THR	2.9
1	B	173	ASP	2.9
1	A	277	HIS	2.8
1	F	294	GLU	2.8
1	G	173	ASP	2.8
1	B	184	ARG	2.8
1	D	294	GLU	2.8
1	D	75	SER	2.7
1	C	173	ASP	2.7
1	A	74	ILE	2.7
1	D	69	LEU	2.7
1	B	77	GLU	2.6
1	G	77	GLU	2.6
1	D	291	THR	2.6
1	A	80	ARG	2.6
1	H	340	GLY	2.6
1	H	54	ARG	2.6
1	A	520	GLN	2.6
1	C	184	ARG	2.5
1	F	277	HIS	2.5
1	C	289	ILE	2.5
1	E	80	ARG	2.5
1	D	71	ASN	2.5
1	A	172	ARG	2.5
1	D	8	LYS	2.5
1	F	172	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	295	ALA	2.4
1	B	15	ARG	2.4
1	F	80	ARG	2.4
1	H	72	ARG	2.4
1	H	33	VAL	2.4
1	E	292	GLU	2.4
1	A	295	ALA	2.4
1	A	8	LYS	2.4
1	H	39	ARG	2.4
1	G	277	HIS	2.4
1	B	196	ARG	2.4
1	D	66	PHE	2.4
1	D	175	THR	2.4
1	A	284	LEU	2.3
1	H	175	THR	2.3
1	A	9	PRO	2.3
1	H	293	PRO	2.3
1	H	520	GLN	2.3
1	A	171	ASP	2.3
1	A	480	ASP	2.3
1	D	56	ALA	2.3
1	H	63	THR	2.3
1	H	71	ASN	2.3
1	D	68	PRO	2.3
1	H	73	LEU	2.3
1	E	480	ASP	2.2
1	G	172	ARG	2.2
1	D	173	ASP	2.2
1	E	173	ASP	2.2
1	C	277	HIS	2.2
1	D	62	LEU	2.2
1	G	289	ILE	2.2
1	E	77	GLU	2.1
1	C	480	ASP	2.1
1	G	495	GLU	2.1
1	D	286	ALA	2.1
1	B	54	ARG	2.1
1	H	43	ARG	2.1
1	B	171	ASP	2.1
1	D	60	TYR	2.1
1	E	9	PRO	2.1
1	C	496	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	340	GLY	2.1
1	C	398	ALA	2.1
1	D	73	LEU	2.1
1	H	68	PRO	2.0
1	D	340	GLY	2.0
1	B	40	ASP	2.0
1	B	291	THR	2.0
1	B	177	LEU	2.0
1	C	285	ASP	2.0
1	G	396	GLY	2.0
1	H	12	GLU	2.0
1	G	74	ILE	2.0
1	H	10	ALA	2.0

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

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
1	MDO	H	149	13/14	0.87	0.11	12,15,21,23	0
1	MDO	B	149	13/14	0.88	0.11	12,14,20,24	0
1	MDO	D	149	13/14	0.89	0.10	14,16,20,21	0
1	MDO	E	149	13/14	0.90	0.11	11,12,16,22	0
1	MDO	C	149	13/14	0.91	0.10	10,11,18,21	0
1	MDO	F	149	13/14	0.91	0.10	9,11,17,23	0
1	MDO	G	149	13/14	0.91	0.10	9,11,15,21	0
1	MDO	A	149	13/14	0.91	0.10	11,13,20,21	0

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.