



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:48 AM GMT

PDB ID : 2O7E
Title : Tyrosine ammonia-lyase from *Rhodobacter sphaeroides* (His89Phe variant), bound to 2-aminoindan-2-phosphonic acid
Authors : Louie, G.V.; Bowman, M.E.; Moffitt, M.C.; Baiga, T.J.; Moore, B.S.; Noel, J.P.
Deposited on : 2006-12-11
Resolution : 1.75 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

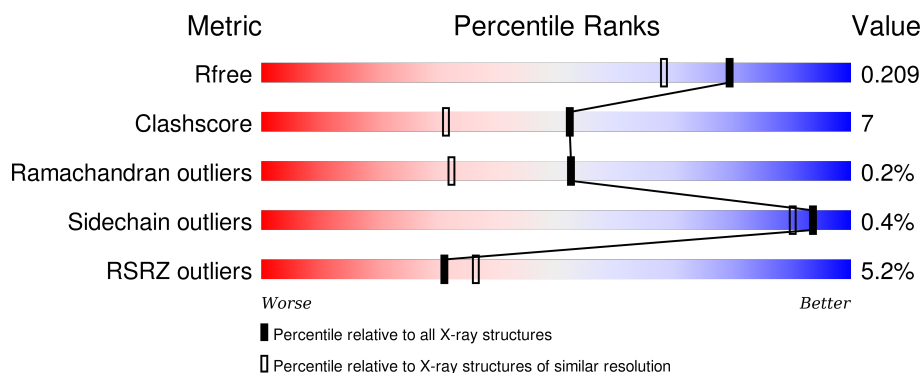
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	91344	1609 (1.76-1.76)
Clashscore	102246	1730 (1.76-1.76)
Ramachandran outliers	100387	1711 (1.76-1.76)
Sidechain outliers	100360	1711 (1.76-1.76)
RSRZ outliers	91569	1610 (1.76-1.76)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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>8%</div> <div>86%</div> <div>13%</div> </div>
1	B	521	<div> <div>5%</div> <div>86%</div> <div>13%</div> </div>
1	C	521	<div> <div>4%</div> <div>83%</div> <div>15%</div> </div>
1	D	521	<div> <div>5%</div> <div>86%</div> <div>13%</div> </div>
1	E	521	<div> <div>4%</div> <div>85%</div> <div>13%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	521	<div><div></div><div>2%87%12%<div>••</div></div></div>
1	G	521	<div><div></div><div>3%87%11%<div>•</div></div></div>
1	H	521	<div><div></div><div>9%85%14%<div>•</div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 33671 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
			3814	2366	729	705	14			
1	B	515	Total	C	N	O	S	0	0	0
			3814	2366	729	705	14			
1	C	514	Total	C	N	O	S	0	0	0
			3807	2361	728	704	14			
1	D	514	Total	C	N	O	S	0	0	0
			3807	2361	728	704	14			
1	E	514	Total	C	N	O	S	0	0	0
			3807	2361	728	704	14			
1	F	514	Total	C	N	O	S	0	0	0
			3807	2361	728	704	14			
1	G	514	Total	C	N	O	S	0	0	0
			3807	2361	728	704	14			
1	H	514	Total	C	N	O	S	0	0	0
			3807	2361	728	704	14			

There are 32 discrepancies between the modelled and reference sequences:

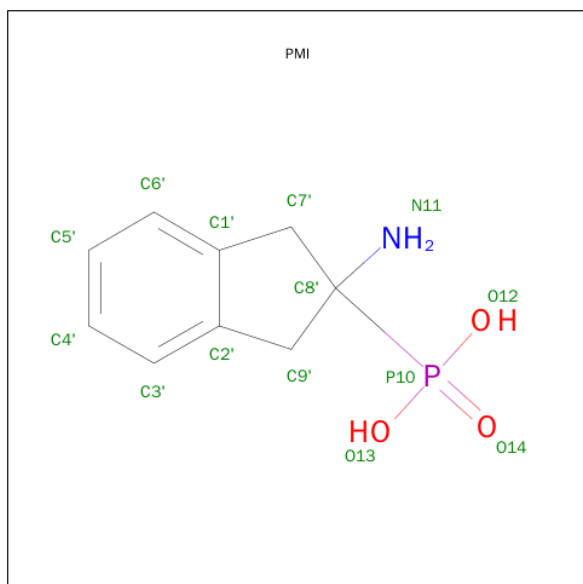
Chain	Residue	Modelled	Actual	Comment	Reference
A	89	PHE	HIS	ENGINEERED	UNP Q3IWB0
A	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
A	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
A	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
B	89	PHE	HIS	ENGINEERED	UNP Q3IWB0
B	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
B	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
B	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
C	89	PHE	HIS	ENGINEERED	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	89	PHE	HIS	ENGINEERED	UNP Q3IWB0

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Chain	Residue	Modelled	Actual	Comment	Reference
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	89	PHE	HIS	ENGINEERED	UNP Q3IWB0
E	149	MDO	ALA	SEE REMARK 999	UNP Q3IWB0
E	149	MDO	SER	SEE REMARK 999	UNP Q3IWB0
E	149	MDO	GLY	SEE REMARK 999	UNP Q3IWB0
F	89	PHE	HIS	ENGINEERED	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	89	PHE	HIS	ENGINEERED	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	89	PHE	HIS	ENGINEERED	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 (2-AMINO-2,3-DIHYDRO-1H-INDEN-2-YL)PHOSPHONIC ACID (three-letter code: PMI) (formula: C₉H₁₂NO₃P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			14	9	1	3	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			14	9	1	3	1		
2	C	1	Total	C	N	O	P	0	0
			14	9	1	3	1		
2	D	1	Total	C	N	O	P	0	0
			14	9	1	3	1		
2	E	1	Total	C	N	O	P	0	0
			14	9	1	3	1		
2	F	1	Total	C	N	O	P	0	0
			14	9	1	3	1		
2	G	1	Total	C	N	O	P	0	0
			14	9	1	3	1		
2	H	1	Total	C	N	O	P	0	0
			14	9	1	3	1		

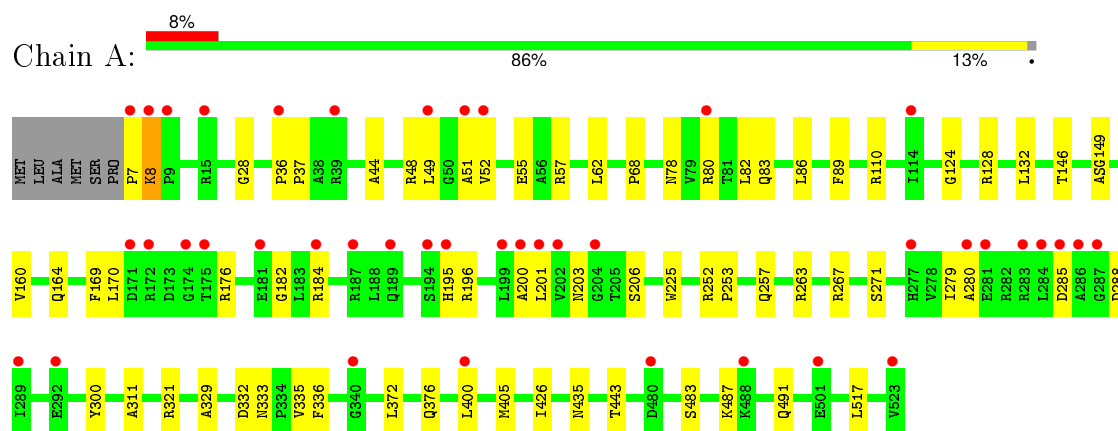
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	325	Total	O	0	0
			325	325		
3	B	373	Total	O	0	0
			373	373		
3	C	414	Total	O	0	0
			414	414		
3	D	355	Total	O	0	0
			355	355		
3	E	380	Total	O	0	0
			380	380		
3	F	479	Total	O	0	0
			479	479		
3	G	413	Total	O	0	0
			413	413		
3	H	350	Total	O	0	0
			350	350		

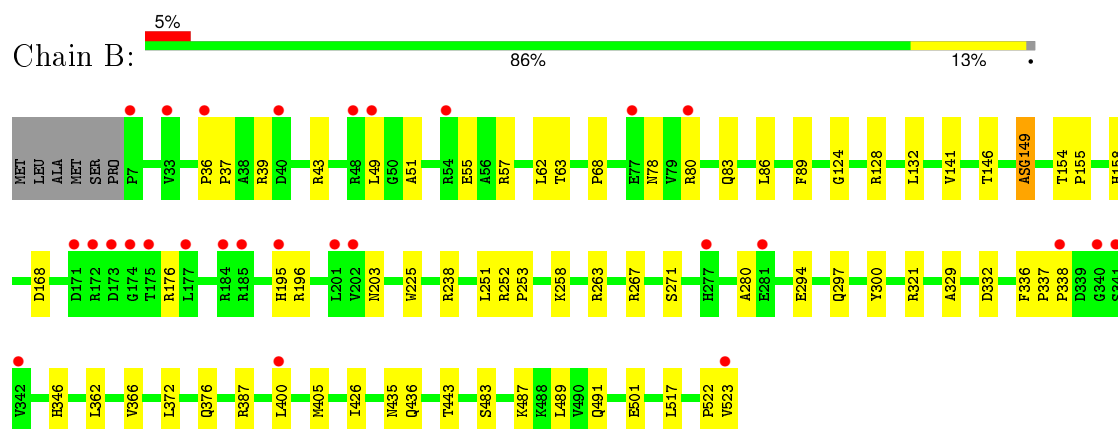
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

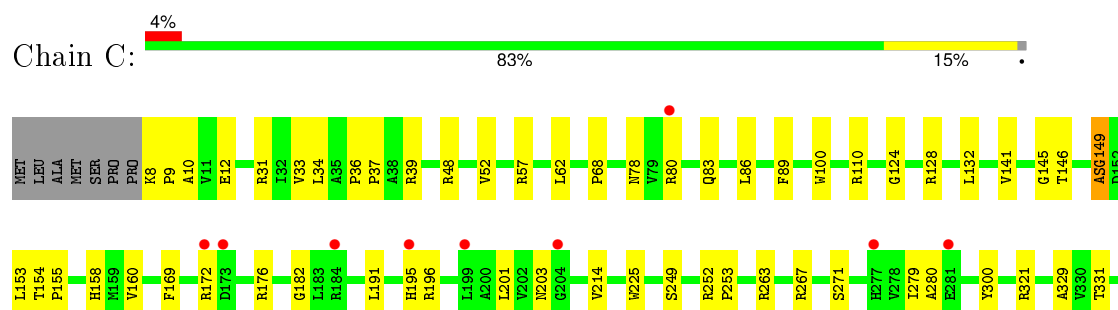
• Molecule 1: Putative histidine ammonia-lyase



• Molecule 1: Putative histidine ammonia-lyase

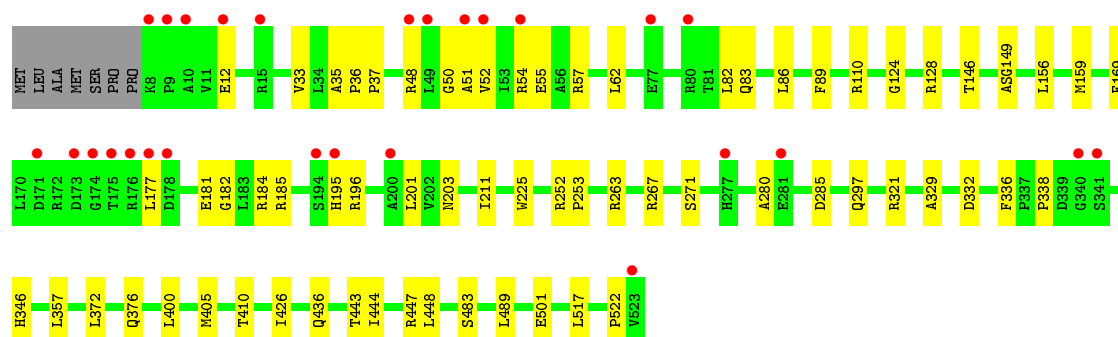
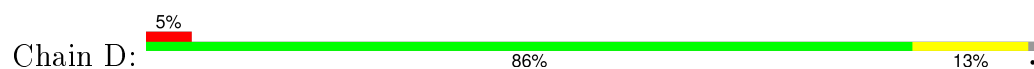


• Molecule 1: Putative histidine ammonia-lyase

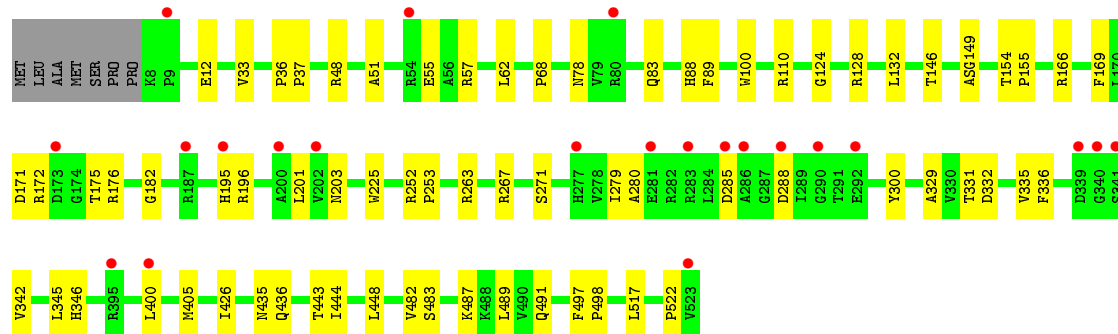
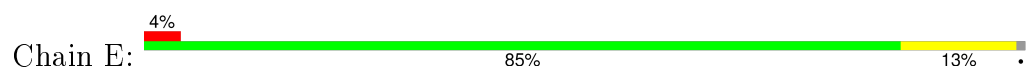




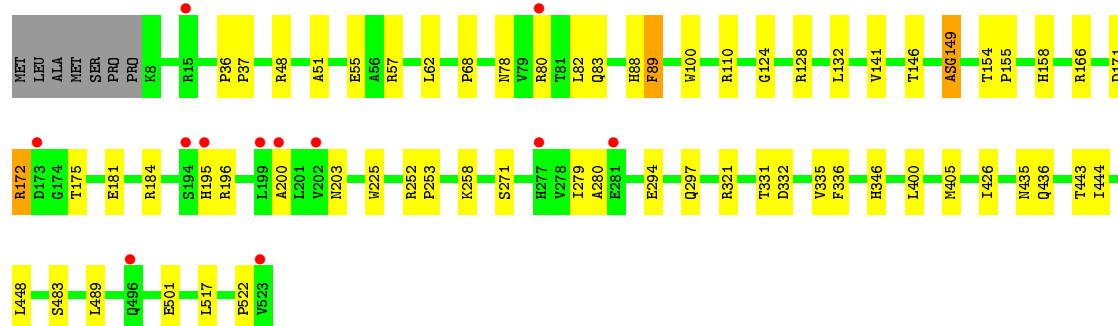
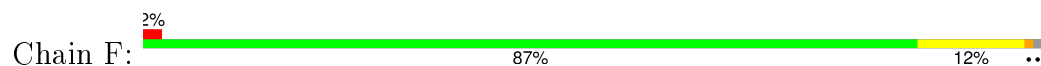
• Molecule 1: Putative histidine ammonia-lyase



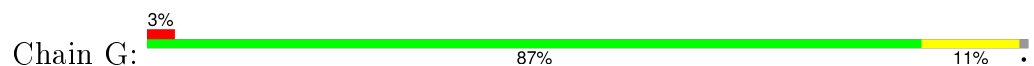
• Molecule 1: Putative histidine ammonia-lyase

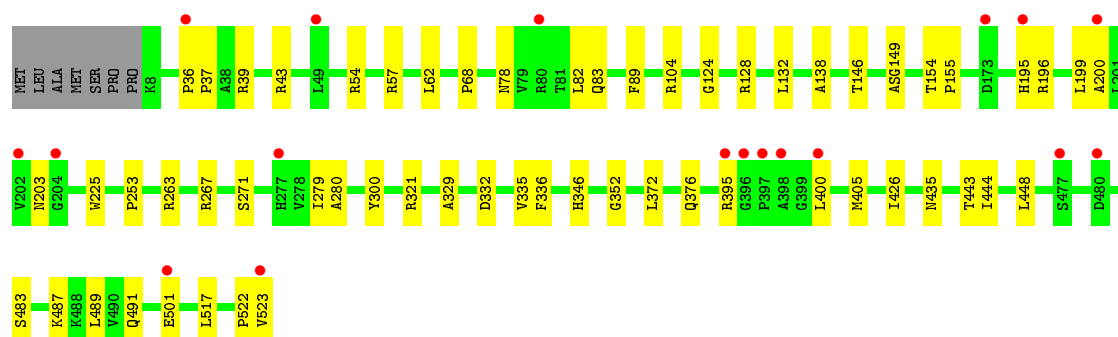


• Molecule 1: Putative histidine ammonia-lyase

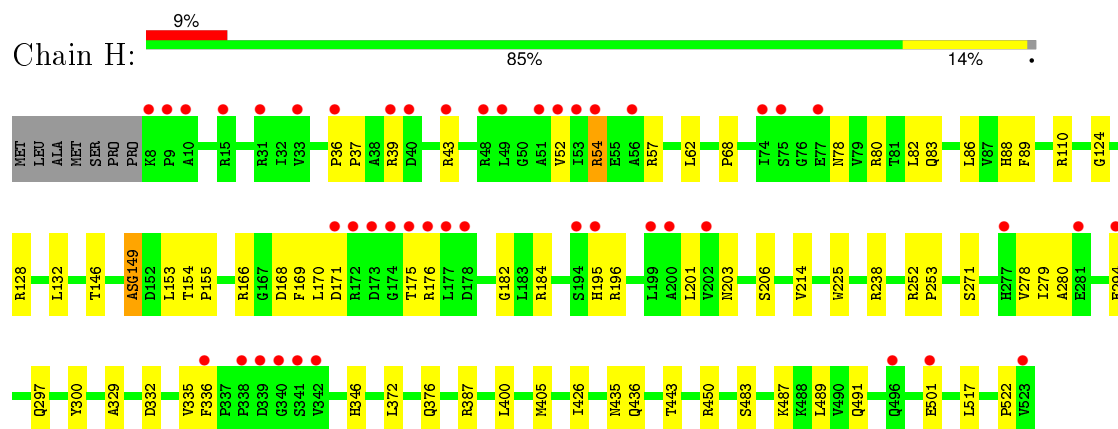


• Molecule 1: Putative histidine ammonia-lyase





• Molecule 1: Putative histidine ammonia-lyase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.57Å 154.84Å 164.53Å 90.00° 94.16° 90.00°	Depositor
Resolution (Å)	500.00 – 1.75 37.68 – 1.75	Depositor EDS
% Data completeness (in resolution range)	96.4 (500.00-1.75) 96.3 (37.68-1.75)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 1.75Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.189 , 0.210 0.189 , 0.209	Depositor DCC
R_{free} test set	21241 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	17.1	Xtriage
Anisotropy	0.243	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 422368 reflections (0.000%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	33671	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% 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, 0.375 respectively for untwinned datasets, and 0.333, 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, PMI

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	0.31	0/3863	0.58	0/5254
1	B	0.31	0/3863	0.58	0/5254
1	C	0.31	0/3855	0.59	0/5243
1	D	0.31	0/3855	0.58	0/5243
1	E	0.31	0/3855	0.58	0/5243
1	F	0.34	0/3855	0.60	0/5243
1	G	0.33	0/3855	0.60	0/5243
1	H	0.30	0/3855	0.57	0/5243
All	All	0.31	0/30856	0.58	0/41966

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.

There are no bond angle outliers.

There are no chirality outliers.

All (16) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	149	MDO	Mainchain,Peptide
1	B	149	MDO	Mainchain,Peptide
1	C	149	MDO	Mainchain,Peptide
1	D	149	MDO	Mainchain,Peptide
1	E	149	MDO	Mainchain,Peptide
1	F	149	MDO	Mainchain,Peptide
1	G	149	MDO	Mainchain,Peptide
1	H	149	MDO	Mainchain,Peptide

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	3814	0	3864	59	0
1	B	3814	0	3865	56	0
1	C	3807	0	3857	64	0
1	D	3807	0	3856	54	0
1	E	3807	0	3857	51	0
1	F	3807	0	3857	57	0
1	G	3807	0	3856	50	0
1	H	3807	0	3857	69	0
2	A	14	0	9	2	0
2	B	14	0	9	3	0
2	C	14	0	9	3	0
2	D	14	0	9	0	0
2	E	14	0	9	2	0
2	F	14	0	9	2	0
2	G	14	0	9	2	0
2	H	14	0	9	3	0
3	A	325	0	0	9	0
3	B	373	0	0	7	0
3	C	414	0	0	12	0
3	D	355	0	0	4	0
3	E	380	0	0	5	0
3	F	479	0	0	11	0
3	G	413	0	0	8	0
3	H	350	0	0	9	0
All	All	33671	0	30941	412	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 7.

All (412) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:83:GLN:HE22	1:H:196:ARG:HA	1.17	1.08
1:A:83:GLN:HE22	1:A:196:ARG:HA	1.21	1.02
1:B:83:GLN:HE22	1:B:196:ARG:HA	1.25	0.97
1:F:200:ALA:HB2	3:F:3012:HOH:O	1.64	0.97
1:G:83:GLN:HE22	1:G:196:ARG:HA	1.28	0.96
1:D:83:GLN:HE22	1:D:196:ARG:HA	1.30	0.94
1:C:83:GLN:HE22	1:C:196:ARG:HA	1.38	0.89
1:F:83:GLN:HE22	1:F:196:ARG:HA	1.41	0.84
1:H:83:GLN:NE2	1:H:196:ARG:HA	1.93	0.82
1:F:62:LEU:HD21	1:F:196:ARG:HB3	1.62	0.81
1:A:62:LEU:HD21	1:A:196:ARG:HB3	1.64	0.79
1:B:51:ALA:O	1:B:55:GLU:HG3	1.83	0.79
1:G:83:GLN:NE2	1:G:196:ARG:HA	1.98	0.78
1:A:200:ALA:HB2	3:A:2945:HOH:O	1.82	0.78
1:B:68:PRO:HG3	1:B:435:ASN:HB2	1.68	0.76
1:A:83:GLN:NE2	1:A:196:ARG:HA	1.99	0.76
1:B:62:LEU:HD21	1:B:196:ARG:HB3	1.68	0.75
1:D:83:GLN:NE2	1:D:196:ARG:HA	2.02	0.74
1:E:196:ARG:HD3	1:E:336:PHE:HE2	1.52	0.74
1:B:196:ARG:HD3	1:B:336:PHE:HE2	1.53	0.74
1:E:83:GLN:HE22	1:E:196:ARG:HA	1.52	0.72
1:G:487:LYS:O	1:G:491:GLN:HG3	1.88	0.72
1:G:62:LEU:HD21	1:G:196:ARG:HB3	1.70	0.72
1:E:487:LYS:O	1:E:491:GLN:HG3	1.90	0.72
1:D:62:LEU:HD21	1:D:196:ARG:HB3	1.70	0.71
1:D:196:ARG:HD3	1:D:336:PHE:HE2	1.54	0.71
1:H:62:LEU:HD21	1:H:196:ARG:HB3	1.71	0.71
1:A:68:PRO:HG3	1:A:435:ASN:HB2	1.72	0.71
1:C:487:LYS:O	1:C:491:GLN:HG3	1.90	0.71
1:D:50:GLY:O	1:D:54:ARG:HG3	1.91	0.70
1:B:83:GLN:NE2	1:B:196:ARG:HA	2.05	0.70
2:E:701:PMI:H5'	1:F:405:MET:HE3	1.74	0.69
1:B:487:LYS:O	1:B:491:GLN:HG3	1.92	0.69
1:F:68:PRO:HG3	1:F:435:ASN:HB2	1.73	0.69
1:B:251:LEU:HD13	1:C:335:VAL:HG21	1.75	0.69
1:C:83:GLN:OE1	1:C:195:HIS:O	2.11	0.68
1:B:258:LYS:HD2	3:B:3514:HOH:O	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:ARG:HD3	1:A:336:PHE:CE2	2.29	0.68
1:H:196:ARG:HD3	1:H:336:PHE:CE2	2.29	0.68
1:E:62:LEU:HD21	1:E:196:ARG:HB3	1.76	0.67
1:A:51:ALA:O	1:A:55:GLU:HG3	1.94	0.67
1:G:200:ALA:HB2	3:G:2988:HOH:O	1.95	0.66
1:D:51:ALA:O	1:D:55:GLU:HG3	1.95	0.66
1:F:196:ARG:HD3	1:F:336:PHE:CE2	2.30	0.66
1:C:196:ARG:HD3	1:C:336:PHE:HE2	1.60	0.66
1:F:83:GLN:NE2	1:F:196:ARG:HA	2.09	0.66
1:E:51:ALA:O	1:E:55:GLU:HG3	1.97	0.65
1:G:68:PRO:HG3	1:G:435:ASN:HB2	1.77	0.65
1:D:196:ARG:HD3	1:D:336:PHE:CE2	2.30	0.65
1:C:62:LEU:HD21	1:C:196:ARG:HB3	1.79	0.65
1:D:297:GLN:HG3	3:D:3086:HOH:O	1.96	0.65
1:H:196:ARG:HD3	1:H:336:PHE:HE2	1.59	0.65
1:E:83:GLN:OE1	1:E:195:HIS:O	2.15	0.64
1:B:196:ARG:HD3	1:B:336:PHE:CE2	2.31	0.64
1:E:196:ARG:HD3	1:E:336:PHE:CE2	2.31	0.64
2:E:701:PMI:H5'	1:F:405:MET:CE	2.28	0.63
1:C:48:ARG:HG2	1:C:48:ARG:HH11	1.63	0.63
1:H:238:ARG:HH11	1:H:387:ARG:NH2	1.96	0.63
2:C:701:PMI:H5'	1:D:405:MET:HE3	1.81	0.63
1:H:214:VAL:HB	1:H:450:ARG:NH2	2.14	0.62
1:F:83:GLN:OE1	1:F:195:HIS:O	2.18	0.62
1:F:171:ASP:OD2	1:F:175:THR:HB	1.99	0.62
1:A:280:ALA:O	1:D:57:ARG:HD2	1.99	0.62
1:A:372:LEU:O	1:A:376:GLN:HG3	2.00	0.62
1:F:124:GLY:O	1:F:128:ARG:HG2	1.99	0.61
1:H:487:LYS:O	1:H:491:GLN:HG3	2.01	0.61
1:F:36:PRO:HB2	1:F:37:PRO:HD3	1.84	0.60
1:H:206:SER:OG	3:H:1354:HOH:O	2.12	0.60
1:B:280:ALA:O	1:C:57:ARG:HD2	2.01	0.60
1:C:83:GLN:NE2	1:C:196:ARG:HA	2.12	0.60
1:C:145:GLY:HA2	1:D:410:THR:HG23	1.84	0.60
1:A:271:SER:HB2	1:A:483:SER:HB3	1.84	0.60
1:A:279:ILE:HD11	1:D:346:HIS:CE1	2.37	0.60
1:C:196:ARG:HD3	1:C:336:PHE:CE2	2.37	0.59
1:E:48:ARG:HG2	1:E:48:ARG:HH11	1.67	0.59
2:A:701:PMI:H5'	1:B:405:MET:HE3	1.84	0.59
1:F:321:ARG:HH21	1:G:321:ARG:HH21	1.50	0.59
1:G:271:SER:HB2	1:G:483:SER:HB3	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:PRO:HB2	1:D:37:PRO:HD3	1.83	0.59
1:G:196:ARG:HD3	1:G:336:PHE:CE2	2.37	0.59
1:C:68:PRO:HG3	1:C:435:ASN:HB2	1.85	0.59
1:B:141:VAL:HG13	1:B:158:HIS:HB2	1.85	0.58
1:F:196:ARG:HD3	1:F:336:PHE:HE2	1.66	0.58
1:H:68:PRO:HG3	1:H:435:ASN:HB2	1.86	0.58
1:D:83:GLN:OE1	1:D:195:HIS:O	2.22	0.58
1:C:124:GLY:O	1:C:128:ARG:HG2	2.03	0.58
1:G:83:GLN:HE22	1:G:196:ARG:CA	2.10	0.57
2:G:701:PMI:H5'	1:H:405:MET:HE3	1.85	0.57
1:G:83:GLN:NE2	1:G:195:HIS:O	2.38	0.56
1:H:271:SER:HB2	1:H:483:SER:HB3	1.86	0.56
1:E:271:SER:HB2	1:E:483:SER:HB3	1.87	0.56
1:C:48:ARG:HG2	1:C:48:ARG:NH1	2.21	0.56
1:E:57:ARG:HD2	1:H:280:ALA:O	2.06	0.56
1:E:280:ALA:O	1:H:57:ARG:HD2	2.06	0.56
1:F:146:THR:HG22	3:F:2088:HOH:O	2.05	0.55
1:G:132:LEU:C	1:G:132:LEU:HD13	2.27	0.55
1:A:78:ASN:HB3	1:B:400:LEU:HD21	1.87	0.55
1:A:225:TRP:CE2	1:A:517:LEU:HD22	2.41	0.55
1:H:39:ARG:HG3	1:H:39:ARG:HH11	1.72	0.55
1:F:82:LEU:HD12	3:F:2155:HOH:O	2.05	0.55
1:H:297:GLN:HG3	3:H:2410:HOH:O	2.06	0.55
1:B:57:ARG:HD2	1:C:280:ALA:O	2.07	0.55
1:H:238:ARG:NH1	1:H:387:ARG:NH2	2.56	0.54
1:F:51:ALA:O	1:F:55:GLU:HG3	2.06	0.54
1:F:280:ALA:O	1:G:57:ARG:HD2	2.06	0.54
1:G:82:LEU:HD12	3:G:1625:HOH:O	2.08	0.54
1:E:400:LEU:HD21	1:F:78:ASN:HB3	1.88	0.54
1:H:335:VAL:HG11	3:H:1982:HOH:O	2.08	0.54
1:D:146:THR:HG22	3:D:2324:HOH:O	2.07	0.54
1:H:83:GLN:OE1	1:H:195:HIS:O	2.25	0.54
1:F:335:VAL:HG11	3:F:1411:HOH:O	2.08	0.54
1:A:57:ARG:HD2	1:D:280:ALA:O	2.08	0.54
1:B:271:SER:HB2	1:B:483:SER:HB3	1.90	0.53
1:D:184:ARG:HG2	1:D:184:ARG:HH11	1.72	0.53
1:B:329:ALA:HB1	1:C:252:ARG:HA	1.89	0.53
1:E:335:VAL:HG11	3:E:1376:HOH:O	2.08	0.53
3:C:2230:HOH:O	1:D:82:LEU:HD12	2.08	0.53
1:B:124:GLY:O	1:B:128:ARG:HG2	2.08	0.53
1:H:83:GLN:HE22	1:H:196:ARG:CA	2.04	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:GLN:OE1	1:A:195:HIS:O	2.26	0.53
1:C:271:SER:HB2	1:C:483:SER:HB3	1.90	0.53
1:H:54:ARG:HG2	1:H:54:ARG:HH11	1.73	0.53
1:D:271:SER:HB2	1:D:483:SER:HB3	1.91	0.53
1:D:169:PHE:CD2	1:D:182:GLY:HA3	2.43	0.53
1:A:196:ARG:HD3	1:A:336:PHE:HE2	1.70	0.53
1:B:436:GLN:HG2	1:C:300:TYR:CZ	2.44	0.53
1:B:68:PRO:HG3	1:B:435:ASN:CB	2.39	0.52
1:F:426:ILE:HD13	3:G:3571:HOH:O	2.09	0.52
1:F:48:ARG:HG2	1:F:48:ARG:HH11	1.73	0.52
1:E:89:PHE:CZ	1:F:405:MET:HE2	2.44	0.52
1:F:225:TRP:CE2	1:F:517:LEU:HD22	2.45	0.52
1:G:78:ASN:HB3	1:H:400:LEU:HD21	1.91	0.52
1:F:141:VAL:HG13	1:F:158:HIS:HB2	1.91	0.52
1:G:225:TRP:CE2	1:G:517:LEU:HD22	2.45	0.52
1:A:405:MET:HE3	2:B:701:PMI:H5'	1.90	0.52
1:E:48:ARG:NH1	1:E:48:ARG:HG2	2.25	0.52
1:E:88:HIS:NE2	1:E:166:ARG:NH2	2.50	0.52
1:B:83:GLN:OE1	1:B:195:HIS:O	2.27	0.51
1:C:176:ARG:HD2	3:C:3285:HOH:O	2.09	0.51
1:G:523:VAL:OXT	1:G:523:VAL:HG12	2.08	0.51
1:G:405:MET:HE3	2:H:701:PMI:H5'	1.90	0.51
1:D:203:ASN:O	1:D:332:ASP:HA	2.10	0.51
1:E:225:TRP:CE2	1:E:517:LEU:HD22	2.46	0.51
1:A:196:ARG:HD3	1:A:336:PHE:CZ	2.46	0.51
1:A:176:ARG:HD2	3:A:3332:HOH:O	2.10	0.51
1:H:225:TRP:CE2	1:H:517:LEU:HD22	2.46	0.51
1:A:487:LYS:O	1:A:491:GLN:HG3	2.11	0.51
1:A:132:LEU:HD13	1:A:132:LEU:C	2.31	0.51
1:H:184:ARG:HG2	1:H:184:ARG:HH11	1.76	0.51
1:E:252:ARG:HA	1:H:329:ALA:HB1	1.92	0.51
1:C:78:ASN:HB3	1:D:400:LEU:HD21	1.93	0.51
1:B:523:VAL:OXT	1:B:523:VAL:HG12	2.11	0.51
1:G:146:THR:HG22	3:G:1967:HOH:O	2.10	0.51
1:F:279:ILE:HD11	1:G:346:HIS:CE1	2.46	0.51
3:A:3436:HOH:O	1:D:426:ILE:HD13	2.11	0.51
1:D:52:VAL:HG13	1:D:57:ARG:HB2	1.93	0.50
1:A:52:VAL:HG13	1:A:57:ARG:HB2	1.93	0.50
1:E:78:ASN:HB3	1:F:400:LEU:HD21	1.94	0.50
1:F:80:ARG:HG2	3:F:2717:HOH:O	2.10	0.50
1:A:426:ILE:HD13	3:D:3398:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:83:GLN:NE2	1:E:196:ARG:HA	2.23	0.50
1:E:68:PRO:HG3	1:E:435:ASN:HB2	1.92	0.50
1:F:271:SER:HB2	1:F:483:SER:HB3	1.93	0.50
1:B:36:PRO:HB2	1:B:37:PRO:HD3	1.92	0.50
1:E:110:ARG:HD3	1:E:201:LEU:HA	1.94	0.50
1:F:57:ARG:HD2	1:G:280:ALA:O	2.12	0.50
1:E:279:ILE:HD11	1:H:346:HIS:CE1	2.46	0.49
1:H:124:GLY:O	1:H:128:ARG:HG2	2.13	0.49
1:A:335:VAL:HG11	3:A:1711:HOH:O	2.12	0.49
2:C:701:PMI:H5'	1:D:405:MET:CE	2.42	0.49
1:H:146:THR:HG22	3:H:2163:HOH:O	2.12	0.49
1:E:154:THR:HB	1:E:155:PRO:CD	2.43	0.49
1:E:346:HIS:CE1	1:H:279:ILE:HD11	2.48	0.49
1:A:329:ALA:HB1	1:D:252:ARG:HA	1.94	0.49
1:E:405:MET:HE3	1:F:89:PHE:CE1	2.47	0.49
1:E:146:THR:HG22	3:E:1932:HOH:O	2.11	0.49
1:C:36:PRO:HB2	1:C:37:PRO:HD3	1.93	0.49
1:E:132:LEU:C	1:E:132:LEU:HD13	2.32	0.49
1:H:83:GLN:NE2	1:H:195:HIS:O	2.46	0.49
1:D:184:ARG:NH1	1:D:184:ARG:HG2	2.27	0.49
1:C:225:TRP:CE2	1:C:517:LEU:HD22	2.47	0.49
1:C:523:VAL:OXT	1:C:523:VAL:HG12	2.13	0.49
1:H:489:LEU:C	1:H:489:LEU:HD23	2.32	0.49
1:B:168:ASP:HB2	1:B:176:ARG:HH21	1.78	0.48
1:H:132:LEU:C	1:H:132:LEU:HD13	2.34	0.48
1:A:86:LEU:HA	1:A:89:PHE:CE2	2.49	0.48
1:B:225:TRP:CE2	1:B:517:LEU:HD22	2.49	0.48
1:B:252:ARG:HA	1:C:329:ALA:HB1	1.96	0.48
1:A:311:ALA:HB1	1:D:357:LEU:HD12	1.95	0.48
1:B:346:HIS:CE1	1:C:279:ILE:HD11	2.48	0.48
1:E:203:ASN:O	1:E:332:ASP:HA	2.14	0.48
1:D:83:GLN:HE22	1:D:196:ARG:CA	2.15	0.48
1:H:52:VAL:HG13	1:H:57:ARG:HB2	1.96	0.48
1:A:321:ARG:HH21	1:D:321:ARG:HH21	1.61	0.48
1:F:68:PRO:HG3	1:F:435:ASN:CB	2.44	0.48
1:G:124:GLY:O	1:G:128:ARG:HG2	2.13	0.48
1:H:86:LEU:HA	1:H:89:PHE:CE2	2.49	0.47
1:A:203:ASN:O	1:A:332:ASP:HA	2.14	0.47
1:H:372:LEU:O	1:H:376:GLN:HG3	2.14	0.47
1:G:395:ARG:CZ	1:G:501:GLU:HG2	2.44	0.47
1:H:171:ASP:OD2	1:H:175:THR:HB	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:80:ARG:HH11	1:H:80:ARG:HG3	1.78	0.47
1:A:405:MET:CE	2:B:701:PMI:H5'	2.44	0.47
1:G:196:ARG:HH21	1:G:199:LEU:HD23	1.80	0.47
1:G:400:LEU:HD21	1:H:78:ASN:HB3	1.97	0.47
1:A:400:LEU:HD21	1:B:78:ASN:HB3	1.95	0.47
1:C:225:TRP:CE2	1:C:522:PRO:HG3	2.50	0.47
1:C:363:ALA:HB1	1:C:452:LYS:HD2	1.97	0.47
1:B:132:LEU:HD13	1:B:132:LEU:C	2.35	0.47
1:A:206:SER:OG	3:A:1478:HOH:O	2.21	0.47
1:C:12:GLU:HA	1:C:33:VAL:HG13	1.96	0.47
1:G:83:GLN:CD	1:G:195:HIS:O	2.53	0.47
1:G:489:LEU:HD23	1:G:489:LEU:C	2.34	0.47
1:C:100:TRP:CZ2	1:C:172:ARG:HG2	2.50	0.47
1:H:294:GLU:OE1	1:H:387:ARG:NH2	2.48	0.47
1:A:169:PHE:CD2	1:A:182:GLY:HA3	2.50	0.47
1:H:36:PRO:N	1:H:37:PRO:HD2	2.29	0.47
1:D:48:ARG:NH2	1:D:338:PRO:O	2.48	0.47
1:H:83:GLN:CD	1:H:195:HIS:O	2.54	0.47
1:D:444:ILE:O	1:D:448:LEU:HG	2.15	0.47
1:B:436:GLN:HG2	1:C:300:TYR:CE2	2.50	0.46
1:C:86:LEU:HA	1:C:89:PHE:CE2	2.50	0.46
1:C:444:ILE:O	1:C:448:LEU:HG	2.14	0.46
1:G:89:PHE:CE1	1:H:405:MET:HE3	2.51	0.46
1:D:124:GLY:O	1:D:128:ARG:HG2	2.14	0.46
1:F:436:GLN:HG2	1:G:300:TYR:CZ	2.50	0.46
1:F:501:GLU:HB2	3:F:2400:HOH:O	2.15	0.46
1:C:154:THR:HB	1:C:155:PRO:CD	2.45	0.46
1:G:39:ARG:HD2	1:G:43:ARG:NH2	2.30	0.46
1:G:335:VAL:HG11	3:G:1135:HOH:O	2.13	0.46
1:G:196:ARG:HD3	1:G:336:PHE:CZ	2.50	0.46
1:H:184:ARG:NH1	1:H:184:ARG:HG2	2.30	0.46
1:C:405:MET:HE3	1:D:89:PHE:CE1	2.51	0.46
1:G:83:GLN:OE1	1:G:195:HIS:O	2.33	0.46
1:G:36:PRO:N	1:G:37:PRO:HD2	2.31	0.46
1:B:86:LEU:HA	1:B:89:PHE:CE2	2.50	0.46
1:B:225:TRP:CE2	1:B:522:PRO:HD3	2.51	0.46
1:A:252:ARG:HA	1:D:329:ALA:HB1	1.98	0.46
1:G:203:ASN:O	1:G:332:ASP:HA	2.15	0.46
1:B:149:MDO:HB21	2:B:701:PMI:H7'1	1.87	0.46
1:H:170:LEU:HD23	1:H:176:ARG:HG2	1.98	0.46
1:E:342:VAL:HG21	1:E:345:LEU:CD2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:88:HIS:NE2	1:F:166:ARG:NH2	2.53	0.46
1:C:34:LEU:HD23	1:C:39:ARG:HG2	1.97	0.46
1:H:39:ARG:O	1:H:43:ARG:HG3	2.15	0.46
1:H:203:ASN:O	1:H:332:ASP:HA	2.16	0.46
1:C:146:THR:HG22	3:C:1832:HOH:O	2.15	0.46
1:D:181:GLU:OE1	1:D:181:GLU:HA	2.16	0.46
1:A:49:LEU:HD21	1:A:196:ARG:H	1.82	0.45
1:H:153:LEU:HB2	3:H:2465:HOH:O	2.15	0.45
1:E:12:GLU:HA	1:E:33:VAL:HG13	1.98	0.45
1:F:489:LEU:HD23	1:F:489:LEU:C	2.37	0.45
1:C:110:ARG:NH2	1:C:331:THR:O	2.49	0.45
1:A:36:PRO:N	1:A:37:PRO:HD2	2.30	0.45
1:A:110:ARG:HD3	1:A:201:LEU:HA	1.98	0.45
1:F:258:LYS:HD2	3:F:5378:HOH:O	2.17	0.45
1:A:83:GLN:HE22	1:A:196:ARG:CA	2.09	0.45
1:H:176:ARG:HD2	3:H:3379:HOH:O	2.16	0.45
1:D:263:ARG:O	1:D:267:ARG:HG2	2.16	0.45
1:B:80:ARG:HB2	3:B:4971:HOH:O	2.17	0.45
1:D:177:LEU:HD21	1:D:185:ARG:NH1	2.31	0.45
1:C:432:ASN:HA	3:C:1165:HOH:O	2.16	0.45
1:F:80:ARG:HB2	3:F:2468:HOH:O	2.15	0.45
1:E:405:MET:CE	1:F:89:PHE:CZ	2.99	0.45
1:G:54:ARG:HD2	3:G:2686:HOH:O	2.17	0.45
1:E:263:ARG:O	1:E:267:ARG:HG2	2.17	0.45
1:H:450:ARG:HG2	1:H:450:ARG:HH11	1.82	0.45
2:A:701:PMI:H5'	1:B:405:MET:CE	2.47	0.45
1:E:110:ARG:NH2	1:E:331:THR:O	2.50	0.45
1:A:7:PRO:O	1:A:8:LYS:C	2.55	0.45
1:F:294:GLU:HG2	3:F:5120:HOH:O	2.16	0.45
1:A:263:ARG:O	1:A:267:ARG:HG2	2.17	0.45
1:H:110:ARG:HD3	1:H:201:LEU:HA	1.99	0.45
1:A:124:GLY:O	1:A:128:ARG:HG2	2.17	0.45
3:E:1223:HOH:O	1:H:278:VAL:HG12	2.16	0.45
1:D:225:TRP:CE2	1:D:517:LEU:HD22	2.52	0.45
1:F:225:TRP:CE2	1:F:522:PRO:HD3	2.51	0.45
1:E:489:LEU:C	1:E:489:LEU:HD23	2.36	0.45
1:D:83:GLN:CD	1:D:195:HIS:O	2.56	0.45
1:B:63:THR:HG22	3:B:4988:HOH:O	2.16	0.45
1:H:82:LEU:HD12	3:H:2439:HOH:O	2.15	0.45
1:G:225:TRP:CE2	1:G:522:PRO:HD3	2.52	0.44
1:A:160:VAL:O	1:A:164:GLN:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:3548:HOH:O	1:C:426:ILE:HD13	2.17	0.44
1:C:110:ARG:HD3	1:C:201:LEU:HA	1.99	0.44
1:G:444:ILE:O	1:G:448:LEU:HG	2.17	0.44
1:H:238:ARG:NH1	1:H:387:ARG:HH22	2.15	0.44
1:C:424:ALA:HB2	1:C:448:LEU:HD12	1.98	0.44
1:A:80:ARG:HB2	3:A:2221:HOH:O	2.16	0.44
1:A:252:ARG:HD3	1:A:257:GLN:OE1	2.16	0.44
1:G:196:ARG:HD3	1:G:336:PHE:HE2	1.80	0.44
1:C:225:TRP:CD2	1:C:522:PRO:HG3	2.53	0.44
1:D:372:LEU:O	1:D:376:GLN:HG3	2.17	0.44
1:F:181:GLU:OE1	1:F:184:ARG:NH1	2.51	0.44
1:G:263:ARG:O	1:G:267:ARG:HG2	2.18	0.44
1:B:154:THR:HB	1:B:155:PRO:CD	2.47	0.44
1:F:154:THR:HB	1:F:155:PRO:CD	2.47	0.44
1:A:196:ARG:HH11	1:A:336:PHE:HE2	1.66	0.44
1:B:489:LEU:HD23	1:B:489:LEU:C	2.38	0.44
1:C:149:MDO:HB21	2:C:701:PMI:H7'1	1.82	0.44
1:A:89:PHE:CE1	1:B:405:MET:HE3	2.52	0.44
2:G:701:PMI:H5'	1:H:405:MET:CE	2.47	0.44
1:H:225:TRP:CE2	1:H:522:PRO:HD3	2.53	0.44
1:B:501:GLU:HB2	3:B:2922:HOH:O	2.18	0.44
1:F:297:GLN:HG3	3:F:2987:HOH:O	2.17	0.44
1:C:263:ARG:O	1:C:267:ARG:HG2	2.17	0.44
1:C:132:LEU:HD13	1:C:132:LEU:C	2.38	0.44
1:H:80:ARG:NH1	1:H:80:ARG:HG3	2.33	0.43
1:E:436:GLN:HG2	1:H:300:TYR:CZ	2.53	0.43
1:C:153:LEU:HD21	3:C:3002:HOH:O	2.17	0.43
1:C:128:ARG:HG3	1:C:191:LEU:HD12	2.00	0.43
1:C:160:VAL:HG13	1:C:191:LEU:HD23	2.00	0.43
1:A:203:ASN:ND2	1:A:333:ASN:HB3	2.33	0.43
1:H:168:ASP:HB2	1:H:176:ARG:HH21	1.83	0.43
1:B:238:ARG:HH21	1:B:387:ARG:HH21	1.64	0.43
1:A:83:GLN:CD	1:A:195:HIS:O	2.57	0.43
1:H:39:ARG:HD3	1:H:43:ARG:HH21	1.83	0.43
1:D:12:GLU:HG2	1:D:35:ALA:HB2	2.01	0.43
1:H:214:VAL:HB	1:H:450:ARG:HH21	1.83	0.43
1:H:149:MDO:HB21	2:H:701:PMI:H7'1	1.85	0.43
1:D:86:LEU:HA	1:D:89:PHE:CE2	2.53	0.43
1:C:249:SER:HB2	3:C:5194:HOH:O	2.18	0.43
1:A:285:ASP:OD1	1:A:288:ASP:OD2	2.37	0.43
1:C:394:HIS:HE1	3:C:2881:HOH:O	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:80:ARG:HG3	1:F:80:ARG:HH11	1.84	0.43
1:D:12:GLU:HA	1:D:33:VAL:HG13	2.00	0.43
1:A:300:TYR:CZ	1:D:436:GLN:HG2	2.53	0.43
1:D:156:LEU:O	1:D:159:MET:HB3	2.19	0.43
1:A:44:ALA:O	1:A:48:ARG:HG3	2.19	0.43
1:B:294:GLU:HG3	1:B:387:ARG:NH2	2.34	0.42
1:E:36:PRO:HB2	1:E:37:PRO:HD3	2.00	0.42
1:F:149:MDO:HB21	2:F:701:PMI:H7'1	1.85	0.42
1:B:426:ILE:HD13	3:C:3493:HOH:O	2.17	0.42
1:F:196:ARG:HD3	1:F:336:PHE:CZ	2.55	0.42
1:E:405:MET:HE3	2:F:701:PMI:H5'	2.00	0.42
1:C:52:VAL:HG23	3:C:1574:HOH:O	2.18	0.42
1:F:346:HIS:CE1	1:G:279:ILE:HD11	2.54	0.42
1:A:184:ARG:HH11	1:A:184:ARG:HG2	1.84	0.42
1:F:252:ARG:HA	1:G:329:ALA:HB1	2.01	0.42
1:G:154:THR:HB	1:G:155:PRO:CD	2.50	0.42
1:F:171:ASP:OD1	1:F:171:ASP:C	2.58	0.42
1:B:372:LEU:O	1:B:376:GLN:HG3	2.19	0.42
1:E:100:TRP:CZ2	1:E:172:ARG:HG2	2.54	0.42
1:E:225:TRP:CE2	1:E:522:PRO:HD3	2.55	0.42
1:B:146:THR:HG22	3:B:2492:HOH:O	2.20	0.42
1:E:171:ASP:OD2	1:E:175:THR:HB	2.19	0.42
1:C:83:GLN:CD	1:C:195:HIS:O	2.57	0.42
3:E:3348:HOH:O	1:H:426:ILE:HD13	2.20	0.42
1:C:12:GLU:HA	1:C:33:VAL:CG1	2.50	0.42
1:E:426:ILE:HD13	3:H:3422:HOH:O	2.19	0.42
1:D:489:LEU:HD23	1:D:489:LEU:C	2.40	0.42
1:E:444:ILE:O	1:E:448:LEU:HG	2.20	0.42
3:F:3660:HOH:O	1:G:426:ILE:HD13	2.20	0.42
1:D:225:TRP:CE2	1:D:522:PRO:HD3	2.55	0.41
1:F:444:ILE:O	1:F:448:LEU:HG	2.19	0.41
1:C:214:VAL:HG13	3:C:4020:HOH:O	2.20	0.41
1:B:321:ARG:HH21	1:C:321:ARG:HH21	1.67	0.41
1:B:39:ARG:HD2	1:B:43:ARG:NH2	2.35	0.41
1:E:169:PHE:CD2	1:E:182:GLY:HA3	2.55	0.41
1:B:49:LEU:HD21	1:B:196:ARG:H	1.85	0.41
1:A:200:ALA:CB	3:A:2945:HOH:O	2.53	0.41
1:B:263:ARG:O	1:B:267:ARG:HG2	2.19	0.41
1:D:83:GLN:NE2	1:D:195:HIS:O	2.53	0.41
1:A:89:PHE:CZ	1:B:405:MET:CE	3.03	0.41
1:B:300:TYR:CZ	1:C:436:GLN:HG2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:169:PHE:CD2	1:C:182:GLY:HA3	2.55	0.41
1:B:297:GLN:HG3	3:B:3097:HOH:O	2.21	0.41
1:C:489:LEU:C	1:C:489:LEU:HD23	2.40	0.41
1:F:83:GLN:HE22	1:F:196:ARG:CA	2.23	0.41
1:G:405:MET:CE	2:H:701:PMI:H5'	2.51	0.41
1:E:497:PHE:HA	1:E:498:PRO:HD3	1.91	0.41
1:E:482:VAL:O	1:E:487:LYS:HE2	2.21	0.41
1:G:352:GLY:HA2	3:G:1417:HOH:O	2.19	0.41
1:F:110:ARG:NH2	1:F:331:THR:O	2.53	0.41
1:F:132:LEU:HD13	1:F:132:LEU:C	2.40	0.41
1:D:211:ILE:HD13	1:D:447:ARG:HG3	2.03	0.41
1:G:200:ALA:CB	3:G:2988:HOH:O	2.60	0.41
1:A:170:LEU:CD2	1:A:176:ARG:HG2	2.51	0.41
1:F:83:GLN:CD	1:F:195:HIS:O	2.58	0.41
1:B:141:VAL:HG13	1:B:158:HIS:CB	2.51	0.41
1:H:171:ASP:OD1	1:H:171:ASP:C	2.59	0.41
1:G:372:LEU:O	1:G:376:GLN:HG3	2.20	0.41
1:H:88:HIS:NE2	1:H:166:ARG:NH2	2.59	0.41
1:C:203:ASN:CG	3:C:3002:HOH:O	2.60	0.41
1:B:337:PRO:HA	1:B:338:PRO:HD3	1.96	0.41
1:E:300:TYR:CZ	1:H:436:GLN:HG2	2.56	0.41
1:B:362:LEU:O	1:B:366:VAL:HG23	2.21	0.41
1:A:82:LEU:HD12	3:A:1805:HOH:O	2.21	0.41
1:D:501:GLU:HB2	3:D:2374:HOH:O	2.20	0.41
1:F:203:ASN:O	1:F:332:ASP:HA	2.21	0.41
1:G:104:ARG:HG3	1:G:138:ALA:HB2	2.03	0.41
1:H:501:GLU:HB2	3:H:3619:HOH:O	2.21	0.41
1:C:141:VAL:HG13	1:C:158:HIS:HB2	2.02	0.41
1:E:124:GLY:O	1:E:128:ARG:HG2	2.21	0.41
1:E:329:ALA:HB1	1:H:252:ARG:HA	2.01	0.41
1:C:195:HIS:CE1	3:C:2333:HOH:O	2.74	0.40
1:A:68:PRO:HG3	1:A:435:ASN:CB	2.46	0.40
1:H:154:THR:HB	1:H:155:PRO:CD	2.51	0.40
1:C:8:LYS:HA	1:C:9:PRO:HD3	1.95	0.40
1:A:8:LYS:HE3	1:A:28:GLY:O	2.21	0.40
1:D:110:ARG:HD3	1:D:201:LEU:HA	2.02	0.40
1:D:285:ASP:C	1:D:285:ASP:OD2	2.60	0.40
1:C:395:ARG:NH1	1:C:501:GLU:HG2	2.35	0.40
1:G:68:PRO:HG3	1:G:435:ASN:CB	2.46	0.40
1:F:100:TRP:CH2	1:F:172:ARG:HD3	2.57	0.40
1:D:51:ALA:HA	1:D:54:ARG:NH1	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:ASN:O	1:B:332:ASP:HA	2.21	0.40
1:E:285:ASP:OD1	1:E:288:ASP:OD2	2.39	0.40
1:H:169:PHE:CD2	1:H:182:GLY:HA3	2.57	0.40
1:C:10:ALA:HB2	1:C:31:ARG:NH2	2.37	0.40
1:E:176:ARG:HD2	3:E:3308:HOH:O	2.21	0.40
1:A:146:THR:HG22	3:A:2274:HOH:O	2.21	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%)	499 (98%)	11 (2%)	2 (0%)	39	19
1	B	512/521 (98%)	504 (98%)	7 (1%)	1 (0%)	52	32
1	C	511/521 (98%)	504 (99%)	6 (1%)	1 (0%)	52	32
1	D	511/521 (98%)	503 (98%)	7 (1%)	1 (0%)	52	32
1	E	511/521 (98%)	502 (98%)	8 (2%)	1 (0%)	52	32
1	F	511/521 (98%)	503 (98%)	7 (1%)	1 (0%)	52	32
1	G	511/521 (98%)	505 (99%)	5 (1%)	1 (0%)	52	32
1	H	511/521 (98%)	499 (98%)	11 (2%)	1 (0%)	52	32
All	All	4090/4168 (98%)	4019 (98%)	62 (2%)	9 (0%)	52	32

All (9) Ramachandran outliers are listed below:

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

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Mol	Chain	Res	Type
1	H	253	PRO
1	B	253	PRO
1	C	253	PRO
1	D	253	PRO
1	G	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%)	383 (100%)	1 (0%)	94	92
1	B	384/389 (99%)	383 (100%)	1 (0%)	94	92
1	C	383/389 (98%)	381 (100%)	2 (0%)	92	87
1	D	383/389 (98%)	382 (100%)	1 (0%)	94	92
1	E	383/389 (98%)	382 (100%)	1 (0%)	94	92
1	F	383/389 (98%)	380 (99%)	3 (1%)	86	77
1	G	383/389 (98%)	382 (100%)	1 (0%)	94	92
1	H	383/389 (98%)	381 (100%)	2 (0%)	92	87
All	All	3066/3112 (98%)	3054 (100%)	12 (0%)	93	90

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	443	THR
1	B	443	THR
1	C	80	ARG
1	C	443	THR
1	D	443	THR
1	E	443	THR
1	F	89	PHE
1	F	172	ARG
1	F	443	THR
1	G	443	THR

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Mol	Chain	Res	Type
1	H	54	ARG
1	H	443	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	GLN
1	A	394	HIS
1	A	491	GLN
1	B	83	GLN
1	B	394	HIS
1	C	83	GLN
1	C	394	HIS
1	D	83	GLN
1	D	189	GLN
1	D	394	HIS
1	D	491	GLN
1	E	83	GLN
1	E	189	GLN
1	E	394	HIS
1	E	516	HIS
1	E	520	GLN
1	F	83	GLN
1	F	394	HIS
1	F	491	GLN
1	G	83	GLN
1	G	394	HIS
1	H	83	GLN
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	A	149	1,2	11,13,14	2.13	3 (27%)	13,18,20	3.67	6 (46%)
1	MDO	B	149	1,2	11,13,14	2.24	3 (27%)	13,18,20	3.55	6 (46%)
1	MDO	C	149	1,2	11,13,14	2.32	3 (27%)	13,18,20	3.48	6 (46%)
1	MDO	D	149	1,2	11,13,14	2.05	2 (18%)	13,18,20	3.37	4 (30%)
1	MDO	E	149	1,2	11,13,14	2.25	3 (27%)	13,18,20	3.52	5 (38%)
1	MDO	F	149	1,2	11,13,14	2.20	3 (27%)	13,18,20	3.44	4 (30%)
1	MDO	G	149	1,2	11,13,14	2.25	3 (27%)	13,18,20	3.59	6 (46%)
1	MDO	H	149	1,2	11,13,14	2.24	3 (27%)	13,18,20	3.58	5 (38%)

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	A	149	1,2	-	0/4/23/24	0/1/1/1
1	MDO	B	149	1,2	-	0/4/23/24	0/1/1/1
1	MDO	C	149	1,2	-	0/4/23/24	0/1/1/1
1	MDO	D	149	1,2	-	0/4/23/24	0/1/1/1
1	MDO	E	149	1,2	-	0/4/23/24	0/1/1/1
1	MDO	F	149	1,2	-	0/4/23/24	0/1/1/1
1	MDO	G	149	1,2	-	0/4/23/24	0/1/1/1
1	MDO	H	149	1,2	-	0/4/23/24	0/1/1/1

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	149	MDO	C2-N3	-4.61	1.30	1.39
1	H	149	MDO	C2-N3	-4.41	1.30	1.39
1	B	149	MDO	C2-N3	-4.26	1.30	1.39
1	G	149	MDO	C2-N3	-4.26	1.30	1.39
1	E	149	MDO	C2-N3	-4.22	1.30	1.39
1	F	149	MDO	C2-N3	-4.15	1.31	1.39
1	A	149	MDO	C2-N3	-4.03	1.31	1.39
1	D	149	MDO	C2-N3	-3.98	1.31	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	149	MDO	CA2-N2	-2.27	1.34	1.39
1	A	149	MDO	CA2-N2	-2.22	1.34	1.39
1	G	149	MDO	CA2-N2	-2.16	1.34	1.39
1	B	149	MDO	CA2-N2	-2.15	1.34	1.39
1	C	149	MDO	CA2-N2	-2.12	1.34	1.39
1	H	149	MDO	CA2-N2	-2.10	1.34	1.39
1	E	149	MDO	CA2-N2	-2.06	1.34	1.39
1	A	149	MDO	O2-C2	4.88	1.33	1.23
1	D	149	MDO	O2-C2	4.89	1.33	1.23
1	G	149	MDO	O2-C2	5.05	1.33	1.23
1	H	149	MDO	O2-C2	5.07	1.33	1.23
1	F	149	MDO	O2-C2	5.12	1.33	1.23
1	B	149	MDO	O2-C2	5.18	1.34	1.23
1	C	149	MDO	O2-C2	5.30	1.34	1.23
1	E	149	MDO	O2-C2	5.35	1.34	1.23

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	MDO	O2-C2-CA2	-5.01	128.24	130.95
1	F	149	MDO	O2-C2-CA2	-4.99	128.25	130.95
1	B	149	MDO	O2-C2-CA2	-4.43	128.55	130.95
1	G	149	MDO	O2-C2-CA2	-4.37	128.58	130.95
1	H	149	MDO	O2-C2-CA2	-4.17	128.69	130.95
1	D	149	MDO	O2-C2-CA2	-3.98	128.80	130.95
1	C	149	MDO	O2-C2-CA2	-3.96	128.81	130.95
1	E	149	MDO	O2-C2-CA2	-3.73	128.93	130.95
1	B	149	MDO	C-CA3-N3	-3.41	105.55	113.00
1	G	149	MDO	C-CA3-N3	-3.34	105.70	113.00
1	E	149	MDO	C-CA3-N3	-3.22	105.96	113.00
1	D	149	MDO	N3-C1-N2	-3.04	109.21	111.56
1	B	149	MDO	N3-C1-N2	-2.94	109.28	111.56
1	C	149	MDO	C-CA3-N3	-2.92	106.61	113.00
1	G	149	MDO	N3-C1-N2	-2.76	109.42	111.56
1	H	149	MDO	C2-CA2-N2	-2.76	106.71	108.91
1	E	149	MDO	N3-C1-N2	-2.74	109.44	111.56
1	C	149	MDO	N3-C1-N2	-2.62	109.53	111.56
1	A	149	MDO	C2-CA2-N2	-2.58	106.85	108.91
1	E	149	MDO	C2-CA2-N2	-2.57	106.86	108.91
1	F	149	MDO	N3-C1-N2	-2.51	109.61	111.56
1	A	149	MDO	N3-C1-N2	-2.51	109.61	111.56
1	D	149	MDO	C2-CA2-N2	-2.46	106.95	108.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	MDO	C-CA3-N3	-2.44	107.66	113.00
1	H	149	MDO	N3-C1-N2	-2.42	109.69	111.56
1	G	149	MDO	C2-CA2-N2	-2.34	107.04	108.91
1	C	149	MDO	C2-CA2-N2	-2.32	107.06	108.91
1	B	149	MDO	C2-CA2-N2	-2.30	107.07	108.91
1	C	149	MDO	CA3-N3-C1	2.02	129.70	127.36
1	B	149	MDO	CA3-N3-C1	2.04	129.72	127.36
1	H	149	MDO	CA3-N3-C1	2.09	129.78	127.36
1	G	149	MDO	CA3-N3-C1	2.09	129.78	127.36
1	A	149	MDO	CA3-N3-C1	2.18	129.89	127.36
1	F	149	MDO	CA3-N3-C1	2.29	130.02	127.36
1	F	149	MDO	CA2-C2-N3	10.09	108.66	103.39
1	D	149	MDO	CA2-C2-N3	10.16	108.69	103.39
1	B	149	MDO	CA2-C2-N3	10.22	108.73	103.39
1	C	149	MDO	CA2-C2-N3	10.37	108.81	103.39
1	E	149	MDO	CA2-C2-N3	10.38	108.81	103.39
1	G	149	MDO	CA2-C2-N3	10.51	108.88	103.39
1	A	149	MDO	CA2-C2-N3	10.82	109.04	103.39
1	H	149	MDO	CA2-C2-N3	10.88	109.07	103.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	149	MDO	1	0
1	C	149	MDO	1	0
1	F	149	MDO	1	0
1	H	149	MDO	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 ligands 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$
2	PMI	A	701	1	14,15,15	1.90	2 (14%)	16,24,24	1.43	2 (12%)
2	PMI	B	701	1	14,15,15	1.90	2 (14%)	16,24,24	1.42	2 (12%)
2	PMI	C	701	1	14,15,15	1.88	4 (28%)	16,24,24	1.41	2 (12%)
2	PMI	D	701	1	14,15,15	1.91	3 (21%)	16,24,24	1.35	1 (6%)
2	PMI	E	701	1	14,15,15	1.84	2 (14%)	16,24,24	1.44	2 (12%)
2	PMI	F	701	1	14,15,15	1.93	3 (21%)	16,24,24	1.41	2 (12%)
2	PMI	G	701	1	14,15,15	1.86	3 (21%)	16,24,24	1.40	2 (12%)
2	PMI	H	701	1	14,15,15	1.88	3 (21%)	16,24,24	1.33	1 (6%)

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
2	PMI	A	701	1	-	0/6/19/19	0/2/2/2
2	PMI	B	701	1	-	0/6/19/19	0/2/2/2
2	PMI	C	701	1	-	0/6/19/19	0/2/2/2
2	PMI	D	701	1	-	0/6/19/19	0/2/2/2
2	PMI	E	701	1	-	0/6/19/19	0/2/2/2
2	PMI	F	701	1	-	0/6/19/19	0/2/2/2
2	PMI	G	701	1	-	0/6/19/19	0/2/2/2
2	PMI	H	701	1	-	0/6/19/19	0/2/2/2

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	701	PMI	P10-O12	-2.07	1.50	1.54
2	C	701	PMI	P10-O12	-2.04	1.50	1.54
2	G	701	PMI	P10-O12	-2.04	1.50	1.54
2	H	701	PMI	C5'-C6'	2.02	1.43	1.38
2	C	701	PMI	C5'-C6'	2.03	1.43	1.38
2	G	701	PMI	C4'-C5'	2.07	1.43	1.38
2	D	701	PMI	C4'-C5'	2.11	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	701	PMI	C5'-C6'	2.15	1.43	1.38
2	C	701	PMI	C4'-C5'	2.16	1.43	1.38
2	H	701	PMI	C4'-C5'	2.18	1.43	1.38
2	E	701	PMI	C4'-C5'	2.19	1.43	1.38
2	B	701	PMI	C4'-C5'	2.23	1.43	1.38
2	A	701	PMI	C4'-C5'	2.25	1.43	1.38
2	F	701	PMI	C4'-C5'	2.30	1.43	1.38
2	G	701	PMI	C1'-C2'	5.10	1.49	1.39
2	H	701	PMI	C1'-C2'	5.18	1.49	1.39
2	D	701	PMI	C1'-C2'	5.21	1.49	1.39
2	E	701	PMI	C1'-C2'	5.21	1.49	1.39
2	C	701	PMI	C1'-C2'	5.22	1.49	1.39
2	F	701	PMI	C1'-C2'	5.31	1.49	1.39
2	B	701	PMI	C1'-C2'	5.36	1.49	1.39
2	A	701	PMI	C1'-C2'	5.39	1.49	1.39

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	701	PMI	C5'-C6'-C1'	-2.16	117.56	120.90
2	C	701	PMI	C5'-C6'-C1'	-2.07	117.70	120.90
2	F	701	PMI	C5'-C6'-C1'	-2.05	117.72	120.90
2	A	701	PMI	C5'-C6'-C1'	-2.04	117.74	120.90
2	B	701	PMI	C5'-C6'-C1'	-2.02	117.77	120.90
2	G	701	PMI	C5'-C6'-C1'	-2.00	117.80	120.90
2	H	701	PMI	C7'-C8'-C9'	4.24	110.02	104.11
2	D	701	PMI	C7'-C8'-C9'	4.35	110.18	104.11
2	C	701	PMI	C7'-C8'-C9'	4.46	110.32	104.11
2	G	701	PMI	C7'-C8'-C9'	4.46	110.33	104.11
2	F	701	PMI	C7'-C8'-C9'	4.48	110.35	104.11
2	E	701	PMI	C7'-C8'-C9'	4.55	110.45	104.11
2	B	701	PMI	C7'-C8'-C9'	4.56	110.47	104.11
2	A	701	PMI	C7'-C8'-C9'	4.60	110.52	104.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	PMI	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	701	PMI	3	0
2	C	701	PMI	3	0
2	E	701	PMI	2	0
2	F	701	PMI	2	0
2	G	701	PMI	2	0
2	H	701	PMI	3	0

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%)	0.46	42 (8%) 14 18	12, 21, 41, 56	0
1	B	514/521 (98%)	0.30	28 (5%) 29 35	9, 18, 37, 52	0
1	C	513/521 (98%)	0.26	19 (3%) 45 51	9, 17, 33, 46	0
1	D	513/521 (98%)	0.35	27 (5%) 30 36	8, 19, 39, 52	0
1	E	513/521 (98%)	0.22	22 (4%) 39 45	9, 17, 36, 62	0
1	F	513/521 (98%)	0.07	12 (2%) 64 70	7, 14, 29, 41	0
1	G	513/521 (98%)	0.18	18 (3%) 48 53	7, 16, 31, 46	0
1	H	513/521 (98%)	0.47	45 (8%) 12 15	8, 19, 40, 55	0
All	All	4106/4168 (98%)	0.29	213 (5%) 31 36	7, 18, 37, 62	0

All (213) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	523	VAL	10.2
1	G	523	VAL	10.2
1	E	523	VAL	9.6
1	H	523	VAL	9.3
1	D	523	VAL	8.7
1	A	7	PRO	8.3
1	B	7	PRO	7.7
1	F	523	VAL	7.6
1	C	523	VAL	7.4
1	A	523	VAL	7.3
1	B	277	HIS	7.3
1	D	277	HIS	7.2
1	A	277	HIS	6.4
1	H	277	HIS	5.6
1	F	277	HIS	5.3
1	H	77	GLU	5.3

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Mol	Chain	Res	Type	RSRZ
1	H	9	PRO	5.1
1	D	9	PRO	5.1
1	D	340	GLY	4.7
1	D	77	GLU	4.6
1	A	285	ASP	4.5
1	A	281	GLU	4.4
1	B	195	HIS	4.4
1	E	283	ARG	4.3
1	A	289	ILE	4.3
1	H	340	GLY	4.2
1	B	175	THR	4.2
1	A	202	VAL	4.1
1	H	173	ASP	4.1
1	H	174	GLY	4.0
1	H	52	VAL	3.9
1	E	277	HIS	3.9
1	A	8	LYS	3.8
1	A	187	ARG	3.8
1	C	340	GLY	3.8
1	B	340	GLY	3.7
1	B	173	ASP	3.7
1	D	195	HIS	3.7
1	B	172	ARG	3.6
1	A	9	PRO	3.6
1	E	285	ASP	3.6
1	C	397	PRO	3.6
1	E	288	ASP	3.5
1	H	336	PHE	3.5
1	A	286	ALA	3.5
1	H	36	PRO	3.5
1	E	281	GLU	3.5
1	G	397	PRO	3.5
1	D	54	ARG	3.4
1	G	202	VAL	3.4
1	B	341	SER	3.4
1	H	338	PRO	3.4
1	D	175	THR	3.3
1	C	277	HIS	3.3
1	H	33	VAL	3.3
1	A	80	ARG	3.3
1	C	172	ARG	3.3
1	D	8	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	G	195	HIS	3.3
1	C	396	GLY	3.2
1	D	52	VAL	3.2
1	G	400	LEU	3.2
1	H	49	LEU	3.2
1	A	184	ARG	3.2
1	E	80	ARG	3.2
1	G	277	HIS	3.2
1	H	195	HIS	3.2
1	H	54	ARG	3.2
1	H	172	ARG	3.2
1	H	202	VAL	3.1
1	F	195	HIS	3.1
1	C	400	LEU	3.1
1	A	49	LEU	3.1
1	D	15	ARG	3.1
1	H	53	ILE	3.1
1	A	195	HIS	3.1
1	E	9	PRO	3.0
1	E	286	ALA	3.0
1	H	48	ARG	3.0
1	A	172	ARG	3.0
1	A	287	GLY	2.9
1	E	340	GLY	2.9
1	B	54	ARG	2.9
1	E	195	HIS	2.9
1	E	202	VAL	2.9
1	H	281	GLU	2.8
1	F	202	VAL	2.8
1	H	39	ARG	2.8
1	D	200	ALA	2.8
1	A	400	LEU	2.8
1	H	199	LEU	2.8
1	C	173	ASP	2.8
1	F	281	GLU	2.7
1	G	200	ALA	2.7
1	A	181	GLU	2.7
1	D	281	GLU	2.7
1	A	52	VAL	2.7
1	H	177	LEU	2.7
1	E	290	GLY	2.7
1	A	36	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	80	ARG	2.7
1	A	200	ALA	2.7
1	E	200	ALA	2.7
1	H	341	SER	2.6
1	D	173	ASP	2.6
1	A	201	LEU	2.6
1	D	341	SER	2.6
1	H	75	SER	2.6
1	H	8	LYS	2.6
1	B	184	ARG	2.6
1	A	283	ARG	2.6
1	H	339	ASP	2.6
1	E	187	ARG	2.6
1	A	199	LEU	2.6
1	B	36	PRO	2.6
1	C	501	GLU	2.6
1	H	342	VAL	2.6
1	H	200	ALA	2.6
1	C	195	HIS	2.6
1	E	400	LEU	2.6
1	H	51	ALA	2.5
1	G	36	PRO	2.5
1	C	199	LEU	2.5
1	D	177	LEU	2.5
1	A	204	GLY	2.5
1	E	341	SER	2.5
1	H	496	GLN	2.5
1	A	15	ARG	2.5
1	E	339	ASP	2.5
1	H	175	THR	2.5
1	F	173	ASP	2.5
1	C	281	GLU	2.5
1	A	39	ARG	2.5
1	B	281	GLU	2.5
1	D	12	GLU	2.5
1	B	342	VAL	2.4
1	A	340	GLY	2.4
1	H	10	ALA	2.4
1	H	56	ALA	2.4
1	B	185	ARG	2.4
1	A	174	GLY	2.4
1	C	341	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	51	ALA	2.4
1	E	173	ASP	2.4
1	G	398	ALA	2.4
1	A	171	ASP	2.4
1	D	194	SER	2.3
1	F	194	SER	2.3
1	A	189	GLN	2.3
1	F	15	ARG	2.3
1	G	80	ARG	2.3
1	A	194	SER	2.3
1	H	194	SER	2.3
1	F	200	ALA	2.3
1	D	80	ARG	2.3
1	D	174	GLY	2.3
1	G	477	SER	2.3
1	B	80	ARG	2.3
1	D	51	ALA	2.3
1	D	176	ARG	2.3
1	E	54	ARG	2.3
1	B	171	ASP	2.3
1	A	280	ALA	2.3
1	B	177	LEU	2.3
1	G	501	GLU	2.3
1	A	480	ASP	2.3
1	H	43	ARG	2.2
1	H	74	ILE	2.2
1	H	171	ASP	2.2
1	G	204	GLY	2.2
1	G	396	GLY	2.2
1	C	80	ARG	2.2
1	A	114	ILE	2.2
1	C	184	ARG	2.2
1	H	178	ASP	2.2
1	A	292	GLU	2.2
1	B	49	LEU	2.2
1	B	201	LEU	2.2
1	H	294	GLU	2.2
1	H	176	ARG	2.2
1	D	48	ARG	2.1
1	A	284	LEU	2.1
1	A	175	THR	2.1
1	D	10	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	501	GLU	2.1
1	D	171	ASP	2.1
1	G	173	ASP	2.1
1	F	496	GLN	2.1
1	B	202	VAL	2.1
1	B	48	ARG	2.1
1	H	15	ARG	2.1
1	B	40	ASP	2.1
1	H	40	ASP	2.1
1	B	400	LEU	2.1
1	G	49	LEU	2.1
1	A	488	LYS	2.1
1	F	199	LEU	2.1
1	H	31	ARG	2.1
1	B	77	GLU	2.1
1	E	292	GLU	2.1
1	G	480	ASP	2.1
1	B	174	GLY	2.0
1	H	501	GLU	2.0
1	D	178	ASP	2.0
1	C	496	GLN	2.0
1	D	49	LEU	2.0
1	B	338	PRO	2.0
1	E	395	ARG	2.0
1	G	395	ARG	2.0
1	B	33	VAL	2.0
1	C	204	GLY	2.0
1	C	339	ASP	2.0
1	C	395	ARG	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
1	MDO	H	149	13/14	0.90	0.15	-	16,19,26,26	0
1	MDO	B	149	13/14	0.95	0.14	-	16,19,23,26	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	MDO	E	149	13/14	0.96	0.11	-	12,14,18,18	0
1	MDO	C	149	13/14	0.96	0.14	-	13,14,18,19	0
1	MDO	F	149	13/14	0.95	0.12	-	12,15,20,21	0
1	MDO	D	149	13/14	0.94	0.12	-	18,20,22,24	0
1	MDO	G	149	13/14	0.94	0.14	-	13,14,17,19	0
1	MDO	A	149	13/14	0.96	0.14	-	15,18,21,21	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	PMI	F	701	14/14	0.92	0.15	1.34	21,25,27,28	0
2	PMI	H	701	14/14	0.90	0.15	0.95	25,31,32,33	0
2	PMI	E	701	14/14	0.92	0.13	0.93	20,24,26,27	0
2	PMI	B	701	14/14	0.91	0.14	0.71	26,29,31,31	0
2	PMI	A	701	14/14	0.92	0.13	0.23	23,26,28,29	0
2	PMI	C	701	14/14	0.93	0.12	0.00	19,22,23,23	0
2	PMI	G	701	14/14	0.95	0.11	-0.39	21,22,23,24	0
2	PMI	D	701	14/14	0.94	0.10	-0.73	26,27,29,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.