



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:27 PM UTC

PDB ID : 7ZB0 / pdb_00007zb0
Title : macrocyclase OphP with 15mer
Authors : Song, H.; Naismith, J.H.
Deposited on : 2022-03-23
Resolution : 2.47 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

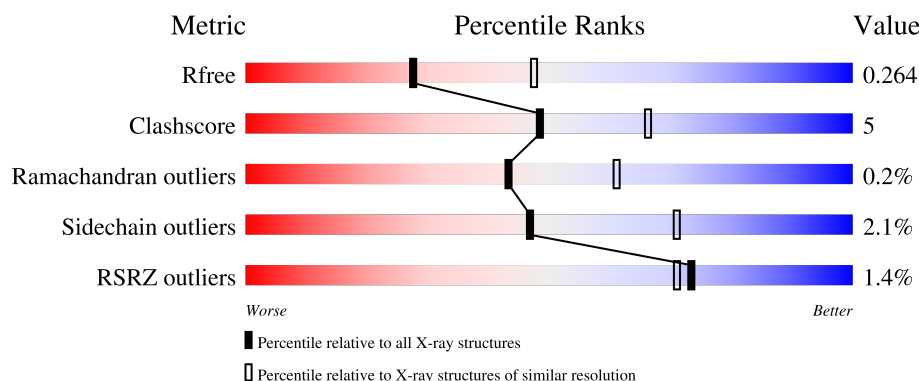
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



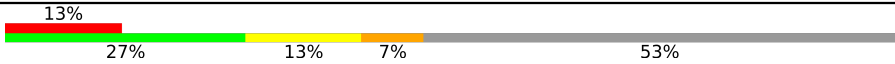
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	745	% 85% 11% ..
1	B	745	% 86% 10% ..
1	C	745	2% 82% 12% 5%
1	D	745	% 85% 11% .
2	E	15	7% 20% 13% 67%

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Mol	Chain	Length	Quality of chain
2	F	15	
2	G	15	
2	H	15	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MVA	E	807	-	-	X	-
2	SAR	G	809	-	-	X	-
2	MVA	H	807	-	-	X	-
2	MVA	H	808	-	-	X	-
2	SAR	H	809	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 23277 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 Prolyl endopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	720	Total	C	N	O	S	0	0	0
			5769	3686	970	1087	26			
1	B	718	Total	C	N	O	S	0	0	0
			5750	3673	968	1083	26			
1	C	706	Total	C	N	O	S	0	0	0
			5667	3627	950	1064	26			
1	D	719	Total	C	N	O	S	0	0	0
			5762	3680	970	1086	26			

- Molecule 2 is a protein called 15mer.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	5	Total	C	N	O	0	0	0
			37	27	5	5			
2	F	7	Total	C	N	O	0	0	0
			64	49	8	7			
2	G	5	Total	C	N	O	0	0	0
			37	27	5	5			
2	H	11	Total	C	N	O	0	0	0
			90	67	12	11			

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is BICARBONATE ION (CCD ID: BCT) (formula: CHO_3).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	1	3		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			7	4	3		
5	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	3	Total 3	Na 3	0	0
6	D	3	Total 3	Na 3	0	0

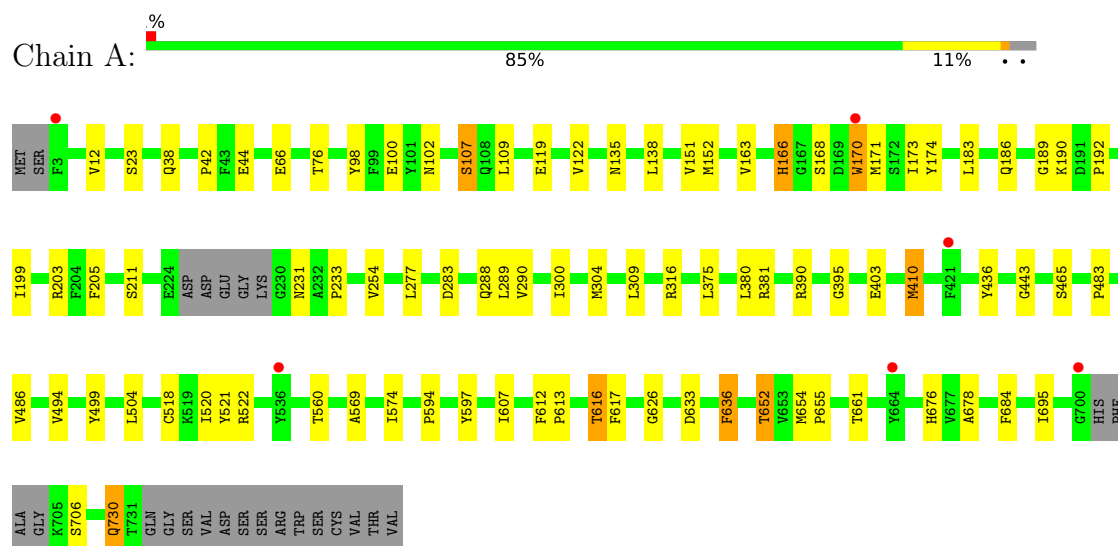
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	7	Total 7	O 7	0	0
7	B	18	Total 18	O 18	0	0
7	C	11	Total 11	O 11	0	0
7	D	17	Total 17	O 17	0	0

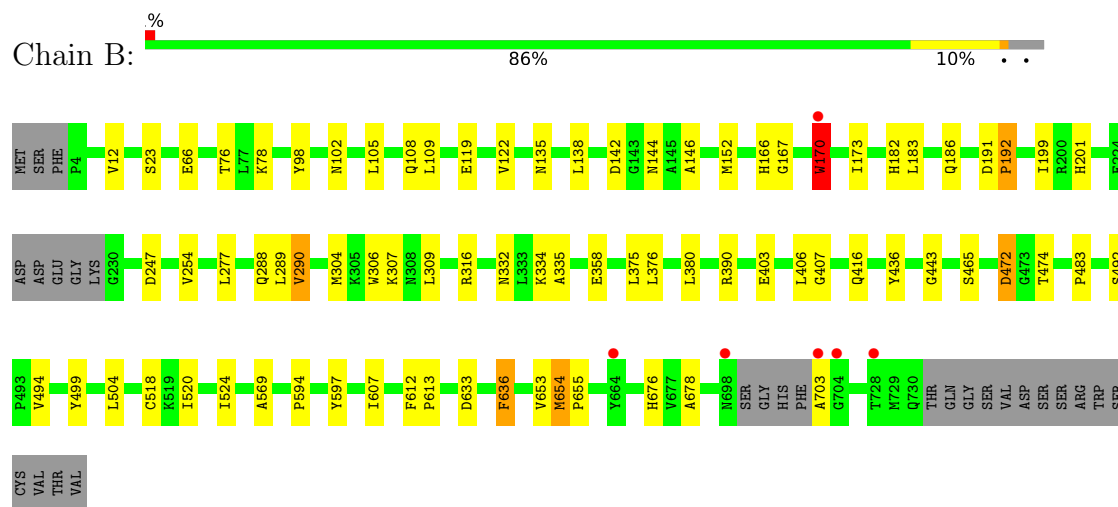
3 Residue-property plots [i](#)

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.

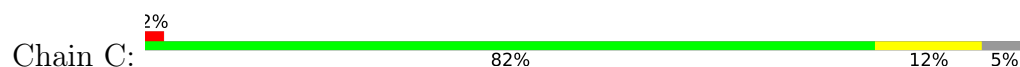
• Molecule 1: Prolyl endopeptidase

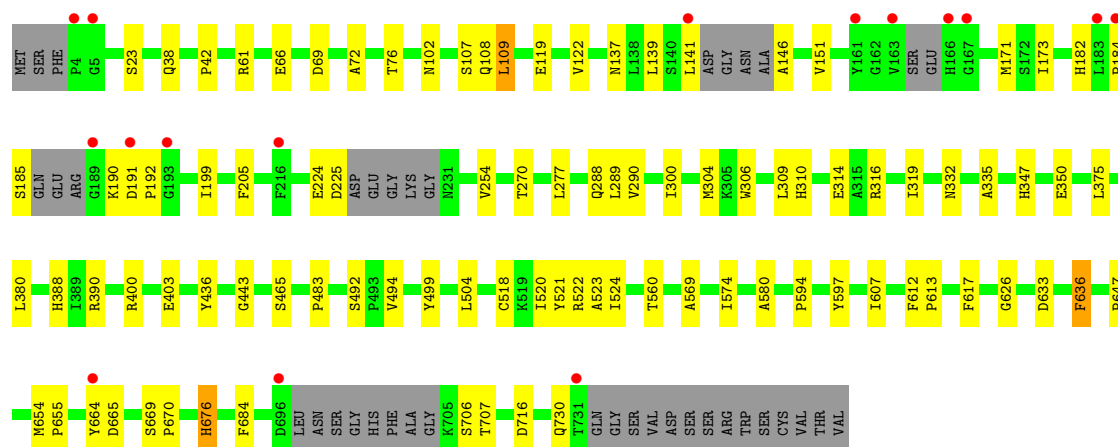


• Molecule 1: Prolyl endopeptidase

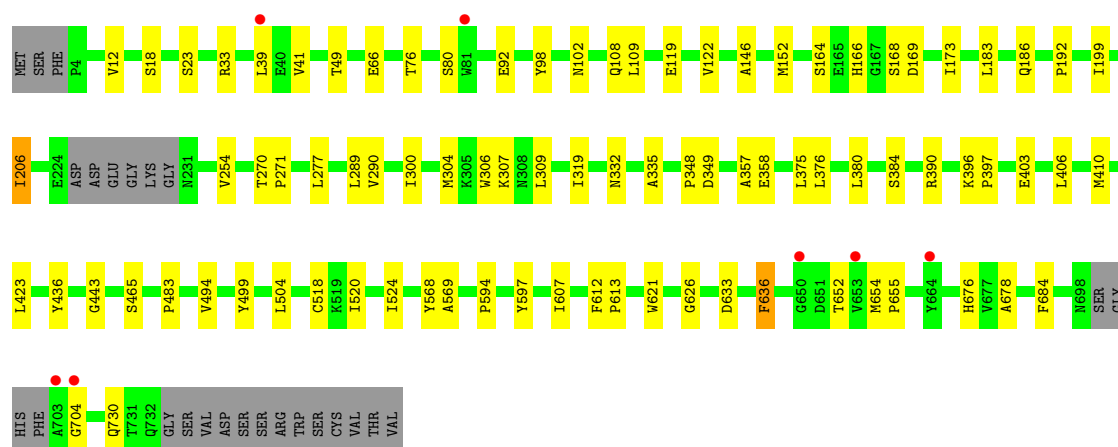
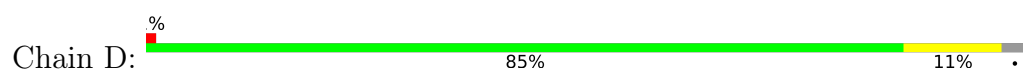


• Molecule 1: Prolyl endopeptidase





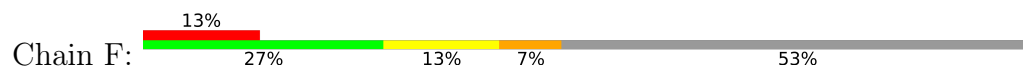
• Molecule 1: Prolyl endopeptidase



• Molecule 2: 15mer

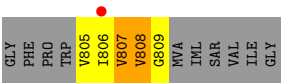


• Molecule 2: 15mer

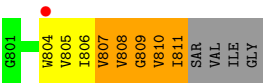


• Molecule 2: 15mer





● Molecule 2: 15mer



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	70.22Å 106.18Å 114.79Å 113.04° 101.67° 93.24°	Depositor
Resolution (Å)	64.35 – 2.47 64.35 – 2.47	Depositor EDS
% Data completeness (in resolution range)	98.6 (64.35-2.47) 98.6 (64.35-2.47)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.229 , 0.266 0.228 , 0.264	Depositor DCC
R_{free} test set	5223 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	55.1	Xtriage
Anisotropy	0.412	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	23277	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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: NA, BCT, SAR, PEG, IML, MVA, EDO

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	1.00	0/5937	1.35	3/8053 (0.0%)
1	B	0.98	0/5917	1.34	3/8025 (0.0%)
1	C	0.98	1/5831 (0.0%)	1.34	7/7906 (0.1%)
1	D	1.01	0/5929	1.34	6/8042 (0.1%)
2	E	0.64	0/7	1.92	0/8
2	F	0.67	0/42	0.87	0/56
2	G	0.15	0/7	1.09	0/8
2	H	0.71	0/46	1.10	0/61
All	All	0.99	1/23716 (0.0%)	1.34	19/32159 (0.1%)

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	1
2	G	0	1
2	H	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	676	HIS	CE1-NE2	5.96	1.38	1.32

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	170	TRP	CA-CB-CG	5.84	124.70	113.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	164	SER	CA-C-N	5.79	128.04	120.28
1	D	164	SER	C-N-CA	5.79	128.04	120.28
1	B	636	PHE	CA-CB-CG	5.69	119.49	113.80
1	D	626	GLY	CA-C-O	-5.58	118.11	122.52
1	A	636	PHE	CA-CB-CG	5.50	119.30	113.80
1	A	626	GLY	CA-C-O	-5.46	118.20	122.52
1	C	665	ASP	CA-C-N	5.39	127.95	120.29
1	C	665	ASP	C-N-CA	5.39	127.95	120.29
1	C	636	PHE	CA-CB-CG	5.37	119.17	113.80
1	B	192	PRO	N-CA-C	-5.32	103.73	111.33
1	A	652	THR	CA-CB-OG1	-5.30	101.64	109.60
1	C	647	PRO	CB-CA-C	-5.23	106.25	111.17
1	D	704	GLY	CA-C-N	5.19	128.22	120.90
1	D	704	GLY	C-N-CA	5.19	128.22	120.90
1	C	626	GLY	CA-C-O	-5.08	118.51	122.52
1	C	42	PRO	CA-C-N	5.00	127.29	120.54
1	C	42	PRO	C-N-CA	5.00	127.29	120.54
1	D	636	PHE	CA-CB-CG	5.00	118.80	113.80

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	730	GLN	Peptide
2	G	808	MVA	Peptide
2	H	809	SAR	Peptide

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	5769	0	5512	53	0
1	B	5750	0	5497	46	0
1	C	5667	0	5420	57	0
1	D	5762	0	5509	59	0
2	E	37	0	50	14	0
2	F	64	0	69	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	37	0	50	10	0
2	H	90	0	101	21	0
3	B	4	0	6	0	0
3	C	8	0	12	0	0
3	D	12	0	18	0	0
4	B	4	0	0	0	0
5	C	7	0	10	0	0
5	D	7	0	10	0	0
6	C	3	0	0	0	0
6	D	3	0	0	0	0
7	A	7	0	0	1	0
7	B	18	0	0	0	0
7	C	11	0	0	0	0
7	D	17	0	0	0	0
All	All	23277	0	22264	249	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 5.

All (249) 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:621:TRP:HE1	2:H:808:MVA:HN2	1.14	1.12
2:F:807:MVA:HG12	2:F:807:MVA:HN2	1.50	0.93
2:H:807:MVA:HN3	2:H:808:MVA:HN1	1.48	0.93
1:D:206:ILE:HD11	2:H:804:TRP:CD2	2.12	0.83
1:D:621:TRP:NE1	2:H:808:MVA:HN2	1.95	0.80
2:H:807:MVA:CN	2:H:808:MVA:HN1	2.11	0.80
1:D:41:VAL:O	1:D:49:THR:HG21	1.84	0.76
1:D:621:TRP:HE1	2:H:808:MVA:CN	1.98	0.73
1:D:277:LEU:HD11	1:D:290:VAL:HG22	1.73	0.71
1:A:100:GLU:OE2	1:A:109:LEU:HD13	1.91	0.70
1:B:277:LEU:HD11	1:B:290:VAL:HG22	1.73	0.69
1:A:277:LEU:HD11	1:A:290:VAL:HG22	1.74	0.69
1:C:277:LEU:HD11	1:C:290:VAL:HG22	1.75	0.69
1:C:171:MET:HE1	1:C:205:PHE:HD1	1.58	0.68
1:A:290:VAL:HG13	1:A:309:LEU:HB3	1.76	0.68
2:E:806:ILE:CG1	2:E:807:MVA:HN1	2.24	0.67
1:D:166:HIS:NE2	1:D:168:SER:HB3	2.08	0.67
2:E:806:ILE:O	2:E:807:MVA:HG12	1.94	0.66
1:C:388:HIS:CE1	1:C:400:ARG:HD3	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:290:VAL:HG13	1:D:309:LEU:HB3	1.77	0.66
1:B:406:LEU:C	1:B:406:LEU:HD23	2.21	0.65
1:A:171:MET:HE1	1:A:205:PHE:HD1	1.61	0.65
1:B:183:LEU:O	1:B:186:GLN:HG3	1.96	0.65
1:B:144:ASN:ND2	1:B:166:HIS:HA	2.12	0.63
1:D:183:LEU:O	1:D:186:GLN:HG3	1.97	0.63
1:A:654:MET:HG3	1:A:684:PHE:CE2	2.33	0.62
1:B:290:VAL:HG13	1:B:309:LEU:HB3	1.81	0.62
1:C:594:PRO:HB3	1:C:654:MET:HE2	1.82	0.62
2:H:807:MVA:N	2:H:808:MVA:HN1	2.15	0.61
1:A:183:LEU:O	1:A:186:GLN:HG3	2.01	0.61
1:A:483:PRO:HD3	1:A:518:CYS:HB3	1.82	0.61
1:A:42:PRO:HG3	1:A:231:ASN:HB2	1.82	0.61
1:C:290:VAL:HG13	1:C:309:LEU:HB3	1.80	0.61
1:C:654:MET:HG3	1:C:684:PHE:CE2	2.36	0.61
1:C:664:TYR:O	1:C:664:TYR:CD1	2.54	0.61
2:H:809:SAR:HN2	2:H:809:SAR:O	2.01	0.61
2:E:806:ILE:HG23	2:E:807:MVA:HN1	1.81	0.60
2:G:805:MVA:HG12	2:G:805:MVA:HN3	1.84	0.60
1:A:617:PHE:HB3	2:E:807:MVA:HG21	1.84	0.60
2:F:807:MVA:HG12	2:F:807:MVA:CN	2.29	0.59
1:C:254:VAL:HG13	1:C:304:MET:HE2	1.85	0.59
1:D:206:ILE:HD11	2:H:804:TRP:CG	2.36	0.59
1:A:594:PRO:HB3	1:A:654:MET:HE2	1.85	0.59
1:C:310:HIS:HE1	1:C:314:GLU:OE1	1.85	0.59
1:B:142:ASP:OD1	1:B:144:ASN:OD1	2.21	0.58
1:D:76:THR:HG21	1:D:520:ILE:HD11	1.85	0.58
1:D:654:MET:HG3	1:D:684:PHE:CE2	2.38	0.58
2:E:805:MVA:HG12	2:E:805:MVA:HN3	1.84	0.58
2:G:808:MVA:HN2	2:G:808:MVA:HG23	1.85	0.58
2:H:808:MVA:HN3	2:H:808:MVA:HG23	1.85	0.57
1:D:33:ARG:HG2	1:D:33:ARG:HH11	1.69	0.57
1:A:163:VAL:CG2	1:A:174:TYR:CE1	2.88	0.57
1:C:171:MET:HE1	1:C:205:PHE:CD1	2.40	0.56
1:A:173:ILE:HB	1:A:199:ILE:HB	1.88	0.56
1:D:483:PRO:HD3	1:D:518:CYS:HB3	1.87	0.56
2:E:806:ILE:CB	2:E:807:MVA:HN1	2.36	0.56
1:C:483:PRO:HD3	1:C:518:CYS:HB3	1.89	0.55
1:B:472:ASP:OD2	1:B:474:THR:OG1	2.19	0.55
1:A:283:ASP:O	2:E:806:ILE:HD13	2.06	0.55
1:C:76:THR:HG21	1:C:520:ILE:HD11	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:806:ILE:HG13	2:G:808:MVA:HG12	1.87	0.55
1:D:594:PRO:HB3	1:D:654:MET:HE2	1.89	0.54
1:A:163:VAL:CG2	1:A:174:TYR:HE1	2.20	0.54
1:D:166:HIS:CE1	1:D:168:SER:HB3	2.42	0.54
1:B:254:VAL:HG13	1:B:304:MET:CE	2.37	0.54
1:C:61:ARG:NH2	1:C:716:ASP:OD1	2.40	0.54
1:D:39:LEU:HA	1:D:49:THR:HG22	1.89	0.54
1:B:483:PRO:HD3	1:B:518:CYS:HB3	1.88	0.54
1:C:182:HIS:HE2	1:C:191:ASP:HB2	1.72	0.54
1:C:347:HIS:HB3	1:C:350:GLU:HG3	1.90	0.54
1:D:173:ILE:HB	1:D:199:ILE:HB	1.90	0.54
2:E:806:ILE:CG1	2:E:807:MVA:CN	2.86	0.54
1:A:171:MET:HE1	1:A:205:PHE:CD1	2.43	0.54
1:C:173:ILE:HB	1:C:199:ILE:HB	1.88	0.54
1:C:580:ALA:CB	2:G:809:SAR:HA2	2.38	0.54
1:B:173:ILE:HB	1:B:199:ILE:HB	1.91	0.53
1:B:182:HIS:HE2	1:B:191:ASP:HB2	1.72	0.53
1:C:119:GLU:O	1:C:122:VAL:HG22	2.08	0.53
1:A:138:LEU:HD13	1:A:186:GLN:OE1	2.09	0.53
1:B:254:VAL:HG13	1:B:304:MET:HE2	1.90	0.53
1:B:375:LEU:HB3	1:B:390:ARG:HB2	1.90	0.53
1:D:76:THR:HG21	1:D:520:ILE:CD1	2.39	0.53
1:D:375:LEU:HB3	1:D:390:ARG:HB2	1.91	0.53
1:A:375:LEU:HB3	1:A:390:ARG:HB2	1.91	0.53
1:B:119:GLU:O	1:B:122:VAL:HG22	2.09	0.53
1:A:254:VAL:HG13	1:A:304:MET:CE	2.39	0.52
1:D:166:HIS:CD2	1:D:168:SER:HB3	2.44	0.52
1:D:633:ASP:HA	1:D:636:PHE:CE2	2.45	0.52
1:B:633:ASP:HA	1:B:636:PHE:CE2	2.45	0.52
2:E:806:ILE:CG2	2:E:807:MVA:HN1	2.40	0.52
1:A:119:GLU:O	1:A:122:VAL:HG22	2.10	0.51
1:C:375:LEU:HB3	1:C:390:ARG:HB2	1.92	0.51
1:C:76:THR:HG21	1:C:520:ILE:CD1	2.40	0.51
2:H:809:SAR:O	2:H:810:MVA:HG12	2.11	0.51
1:A:254:VAL:HG13	1:A:304:MET:HE2	1.92	0.51
1:A:494:VAL:CG2	1:A:569:ALA:HB2	2.41	0.51
1:A:633:ASP:HA	1:A:636:PHE:CE2	2.46	0.51
1:D:307:LYS:HG2	1:D:348:PRO:CB	2.41	0.51
2:H:808:MVA:O	2:H:808:MVA:HG13	2.11	0.50
1:B:167:GLY:O	1:B:703:ALA:HB2	2.11	0.50
1:D:102:ASN:HB2	1:D:109:LEU:HD23	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:349:ASP:OD1	1:D:349:ASP:N	2.42	0.50
1:C:633:ASP:HA	1:C:636:PHE:CE2	2.47	0.50
1:B:102:ASN:HB2	1:B:109:LEU:HD23	1.93	0.50
1:D:607:ILE:HG21	1:D:676:HIS:CG	2.47	0.50
1:A:12:VAL:HG21	1:A:678:ALA:HB1	1.93	0.50
1:B:654:MET:HG3	1:B:655:PRO:HD2	1.93	0.50
1:C:254:VAL:HG13	1:C:304:MET:CE	2.42	0.50
1:D:119:GLU:O	1:D:122:VAL:HG22	2.11	0.50
1:D:12:VAL:HG21	1:D:678:ALA:HB1	1.94	0.49
1:C:597:TYR:O	1:C:655:PRO:HB3	2.13	0.49
1:D:384:SER:OG	1:D:406:LEU:HD12	2.12	0.49
1:D:524:ILE:CD1	1:D:568:TYR:HB3	2.43	0.49
2:G:808:MVA:HN2	2:G:808:MVA:CG2	2.42	0.49
2:H:807:MVA:HN2	2:H:807:MVA:HG22	1.93	0.49
2:H:809:SAR:O	2:H:809:SAR:CN	2.60	0.49
1:B:406:LEU:C	1:B:406:LEU:CD2	2.86	0.49
1:A:607:ILE:HG21	1:A:676:HIS:CG	2.47	0.49
1:B:12:VAL:HG21	1:B:678:ALA:HB1	1.94	0.49
1:C:607:ILE:HG21	1:C:676:HIS:CG	2.48	0.49
1:D:494:VAL:CG2	1:D:569:ALA:HB2	2.43	0.49
1:B:494:VAL:CG2	1:B:569:ALA:HB2	2.42	0.49
1:C:664:TYR:O	1:C:664:TYR:CG	2.66	0.48
1:C:664:TYR:CD1	1:C:664:TYR:C	2.90	0.48
1:A:189:GLY:O	1:A:190:LYS:HG3	2.14	0.48
1:A:289:LEU:HD23	1:A:289:LEU:C	2.38	0.48
1:A:594:PRO:HB3	1:A:654:MET:CE	2.43	0.48
1:A:107:SER:HB3	1:A:706:SER:HB2	1.96	0.48
1:D:254:VAL:HG13	1:D:304:MET:HE2	1.95	0.48
1:B:472:ASP:O	1:B:472:ASP:OD1	2.31	0.48
1:B:607:ILE:HG21	1:B:676:HIS:CG	2.49	0.48
1:C:69:ASP:HA	1:C:72:ALA:HB3	1.96	0.48
1:D:304:MET:HE3	1:D:306:TRP:CZ2	2.49	0.48
1:A:170:TRP:HB3	1:A:203:ARG:NH2	2.29	0.47
1:D:18:SER:O	1:D:33:ARG:HD3	2.14	0.47
1:B:597:TYR:O	1:B:655:PRO:HB3	2.14	0.47
1:C:107:SER:HB2	1:C:137:ASN:OD1	2.14	0.47
1:C:580:ALA:HB1	2:G:809:SAR:HA2	1.96	0.47
1:D:410:MET:HE2	1:D:423:LEU:HD23	1.97	0.47
1:A:163:VAL:HG22	1:A:174:TYR:CE1	2.49	0.47
1:A:597:TYR:O	1:A:655:PRO:HB3	2.15	0.47
1:D:597:TYR:O	1:D:655:PRO:HB3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:289:LEU:C	1:B:289:LEU:HD23	2.40	0.47
1:D:254:VAL:HG13	1:D:304:MET:CE	2.45	0.47
1:C:580:ALA:CB	2:G:809:SAR:C	2.94	0.46
1:C:594:PRO:HB3	1:C:654:MET:CE	2.44	0.46
1:D:289:LEU:HD23	1:D:290:VAL:N	2.30	0.46
2:H:810:MVA:HA	2:H:811:IML:HN1	1.80	0.46
1:D:652:THR:HG21	1:D:654:MET:HE3	1.98	0.46
2:E:806:ILE:HG13	2:E:807:MVA:CN	2.45	0.46
1:B:76:THR:HG21	1:B:520:ILE:HD11	1.97	0.46
1:D:307:LYS:HG2	1:D:348:PRO:HB3	1.97	0.46
2:H:808:MVA:HN3	2:H:808:MVA:CG2	2.41	0.46
1:C:520:ILE:O	1:C:730:GLN:HG3	2.16	0.45
1:A:135:ASN:HB3	1:A:138:LEU:HD12	1.98	0.45
1:B:76:THR:HG21	1:B:520:ILE:CD1	2.47	0.45
1:C:499:TYR:CD2	1:C:504:LEU:HD23	2.52	0.45
1:D:289:LEU:HD23	1:D:289:LEU:C	2.41	0.45
1:C:580:ALA:HB3	2:G:809:SAR:C	2.46	0.45
1:C:494:VAL:CG2	1:C:569:ALA:HB2	2.47	0.45
2:E:806:ILE:HG12	2:E:807:MVA:HN1	1.95	0.45
1:B:289:LEU:HD23	1:B:290:VAL:N	2.32	0.45
1:D:98:TYR:CE1	1:D:152:MET:HE1	2.52	0.45
1:B:612:PHE:N	1:B:613:PRO:CD	2.80	0.44
1:D:206:ILE:CD1	2:H:804:TRP:CG	2.99	0.44
1:A:410:MET:HE2	1:A:410:MET:HB2	1.93	0.44
1:A:76:THR:HG21	1:A:520:ILE:HD11	2.00	0.44
1:B:406:LEU:HD23	1:B:407:GLY:N	2.32	0.44
1:D:436:TYR:CZ	1:D:443:GLY:HA3	2.52	0.44
1:C:289:LEU:C	1:C:289:LEU:HD23	2.42	0.44
1:B:499:TYR:CD2	1:B:504:LEU:HD23	2.52	0.44
1:A:499:TYR:CD2	1:A:504:LEU:HD23	2.53	0.44
1:B:594:PRO:HB3	1:B:654:MET:CE	2.47	0.44
1:D:520:ILE:O	1:D:730:GLN:HG3	2.18	0.44
1:C:436:TYR:CZ	1:C:443:GLY:HA3	2.53	0.43
2:E:808:MVA:HA	2:E:809:SAR:HN1	1.65	0.43
1:B:436:TYR:CZ	1:B:443:GLY:HA3	2.53	0.43
2:H:804:TRP:HA	2:H:805:MVA:HA	1.79	0.43
1:A:289:LEU:HD23	1:A:290:VAL:N	2.34	0.43
1:B:135:ASN:HB3	1:B:138:LEU:HD12	1.99	0.43
1:C:612:PHE:N	1:C:613:PRO:CD	2.81	0.43
1:A:233:PRO:HA	1:A:616:THR:O	2.18	0.43
1:A:381:ARG:NH2	7:A:802:HOH:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:332:ASN:HA	1:B:335:ALA:O	2.17	0.43
1:C:107:SER:HB3	1:C:706:SER:HB2	2.00	0.43
1:C:184:PRO:O	1:C:185:SER:C	2.61	0.43
1:C:523:ALA:C	1:C:524:ILE:HD12	2.44	0.43
1:D:396:LYS:HG3	1:D:397:PRO:HD2	2.01	0.43
1:A:612:PHE:N	1:A:613:PRO:CD	2.82	0.43
1:B:304:MET:HE3	1:B:306:TRP:CZ2	2.54	0.43
1:C:102:ASN:HD22	1:C:109:LEU:HD12	1.83	0.43
1:D:612:PHE:CG	1:D:613:PRO:HD3	2.54	0.43
2:E:808:MVA:O	2:E:808:MVA:HG12	2.19	0.43
1:B:108:GLN:HE22	1:B:146:ALA:HA	1.84	0.42
1:D:594:PRO:HB3	1:D:654:MET:CE	2.48	0.42
1:B:182:HIS:HE2	1:B:191:ASP:CB	2.31	0.42
1:B:334:LYS:HG2	1:D:357:ALA:HB1	2.02	0.42
2:H:806:ILE:HA	2:H:807:MVA:HN1	1.72	0.42
1:B:98:TYR:CE1	1:B:152:MET:HE1	2.55	0.42
1:C:107:SER:O	1:C:707:THR:HG23	2.19	0.42
1:C:612:PHE:CG	1:C:613:PRO:HD3	2.55	0.42
1:D:206:ILE:CD1	2:H:804:TRP:CD2	2.92	0.42
1:D:332:ASN:HA	1:D:335:ALA:O	2.19	0.42
1:C:332:ASN:HA	1:C:335:ALA:O	2.18	0.42
1:D:335:ALA:HB1	1:D:358:GLU:HB2	2.02	0.42
1:A:436:TYR:CZ	1:A:443:GLY:HA3	2.55	0.42
1:B:186:GLN:NE2	1:B:192:PRO:HG2	2.35	0.42
1:B:612:PHE:CG	1:B:613:PRO:HD3	2.55	0.42
2:H:807:MVA:HG13	2:H:809:SAR:HN1	2.02	0.42
1:A:98:TYR:CE1	1:A:152:MET:HE1	2.55	0.42
1:A:102:ASN:HB2	1:A:109:LEU:HD23	2.01	0.42
1:C:289:LEU:HD23	1:C:290:VAL:N	2.34	0.42
1:D:499:TYR:CD2	1:D:504:LEU:HD23	2.54	0.42
1:D:612:PHE:N	1:D:613:PRO:CD	2.82	0.42
1:D:376:LEU:C	1:D:376:LEU:HD23	2.45	0.41
2:F:807:MVA:HA	2:F:808:MVA:HN1	1.77	0.41
1:B:288:GLN:CD	1:B:316:ARG:HG3	2.45	0.41
1:A:521:TYR:O	1:A:522:ARG:C	2.63	0.41
1:A:612:PHE:CG	1:A:613:PRO:HD3	2.55	0.41
1:D:92:GLU:HG3	1:D:98:TYR:CD1	2.55	0.41
1:B:170:TRP:HZ2	1:B:201:HIS:CD2	2.38	0.41
1:C:139:LEU:CD2	1:C:191:ASP:HB3	2.50	0.41
1:C:224:GLU:O	1:C:225:ASP:C	2.63	0.41
1:D:108:GLN:HE22	1:D:146:ALA:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:807:MVA:HA	2:E:808:MVA:HN1	1.68	0.41
2:G:807:MVA:HA	2:G:808:MVA:HN1	1.74	0.41
1:A:168:SER:HB3	1:A:170:TRP:CE2	2.55	0.41
1:A:171:MET:HE1	1:A:205:PHE:HB2	2.03	0.41
1:A:483:PRO:HB2	1:A:486:VAL:HG13	2.03	0.41
1:C:108:GLN:HE22	1:C:146:ALA:HA	1.86	0.41
1:C:669:SER:HA	1:C:670:PRO:HD3	1.96	0.41
1:D:166:HIS:NE2	1:D:168:SER:CB	2.81	0.41
1:A:520:ILE:O	1:A:730:GLN:HG3	2.22	0.40
1:B:376:LEU:HD23	1:B:376:LEU:C	2.46	0.40
1:D:270:THR:HA	1:D:319:ILE:HG21	2.02	0.40
2:F:806:ILE:HA	2:F:807:MVA:HN1	1.76	0.40
1:A:288:GLN:CD	1:A:316:ARG:HG3	2.46	0.40
1:A:390:ARG:NH1	1:A:395:GLY:O	2.53	0.40
1:A:494:VAL:HG23	1:A:569:ALA:CB	2.52	0.40
1:A:560:THR:HG23	1:A:574:ILE:HG21	2.03	0.40
1:A:661:THR:HG23	1:A:695:ILE:HD13	2.03	0.40
1:C:288:GLN:CD	1:C:316:ARG:HG3	2.47	0.40
1:C:560:THR:HG23	1:C:574:ILE:HG21	2.02	0.40
1:C:182:HIS:HE2	1:C:191:ASP:CB	2.32	0.40
1:C:304:MET:HE3	1:C:306:TRP:CZ2	2.56	0.40
1:C:521:TYR:O	1:C:522:ARG:C	2.64	0.40
1:B:335:ALA:HB1	1:B:358:GLU:HB2	2.03	0.40
1:C:270:THR:HA	1:C:319:ILE:HG21	2.03	0.40
1:C:617:PHE:CD1	2:G:807:MVA:HG21	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	714/745 (96%)	673 (94%)	38 (5%)	3 (0%)	30 46

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	712/745 (96%)	674 (95%)	37 (5%)	1 (0%)	48	66
1	C	694/745 (93%)	652 (94%)	41 (6%)	1 (0%)	48	66
1	D	713/745 (96%)	673 (94%)	38 (5%)	2 (0%)	36	53
2	E	1/15 (7%)	0	1 (100%)	0	100	100
2	F	3/15 (20%)	3 (100%)	0	0	100	100
2	G	1/15 (7%)	1 (100%)	0	0	100	100
2	H	4/15 (27%)	4 (100%)	0	0	100	100
All	All	2842/3040 (94%)	2680 (94%)	155 (6%)	7 (0%)	43	61

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	HIS
1	A	192	PRO
1	C	192	PRO
1	D	192	PRO
1	D	169	ASP
1	B	653	VAL
1	A	616	THR

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	623/644 (97%)	608 (98%)	15 (2%)	43	67
1	B	620/644 (96%)	604 (97%)	16 (3%)	40	65
1	C	613/644 (95%)	601 (98%)	12 (2%)	48	72
1	D	622/644 (97%)	613 (99%)	9 (1%)	59	79
2	E	1/6 (17%)	0	1 (100%)	0	0
2	F	4/6 (67%)	4 (100%)	0	100	100
2	G	1/6 (17%)	1 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	4/6 (67%)	4 (100%)	0	100	100
All	All	2488/2600 (96%)	2435 (98%)	53 (2%)	47	71

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	SER
1	A	38	GLN
1	A	44	GLU
1	A	66	GLU
1	A	107	SER
1	A	151	VAL
1	A	166	HIS
1	A	170	TRP
1	A	211	SER
1	A	300	ILE
1	A	380	LEU
1	A	403	GLU
1	A	410	MET
1	A	465	SER
1	A	652	THR
1	B	23	SER
1	B	66	GLU
1	B	78	LYS
1	B	105	LEU
1	B	170	TRP
1	B	247	ASP
1	B	290	VAL
1	B	307	LYS
1	B	380	LEU
1	B	403	GLU
1	B	416	GLN
1	B	465	SER
1	B	472	ASP
1	B	492	SER
1	B	524	ILE
1	B	654	MET
1	C	23	SER
1	C	38	GLN
1	C	66	GLU
1	C	109	LEU
1	C	141	LEU

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Mol	Chain	Res	Type
1	C	151	VAL
1	C	190	LYS
1	C	300	ILE
1	C	380	LEU
1	C	403	GLU
1	C	465	SER
1	C	492	SER
1	D	23	SER
1	D	66	GLU
1	D	80	SER
1	D	206	ILE
1	D	271	PRO
1	D	300	ILE
1	D	380	LEU
1	D	403	GLU
1	D	465	SER
2	E	806	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	268	GLN
1	A	550	GLN
1	A	592	GLN
1	B	38	GLN
1	B	201	HIS
1	B	303	ASN
1	B	308	ASN
1	B	550	GLN
1	B	592	GLN
1	C	38	GLN
1	C	310	HIS
1	C	550	GLN
1	C	592	GLN
1	D	550	GLN
1	D	592	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

17 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
2	MVA	F	807	2	6,7,8	0.68	0	6,8,10	1.63	2 (33%)
2	MVA	G	807	2	6,7,8	0.65	0	6,8,10	2.33	3 (50%)
2	SAR	G	809	2	3,4,5	0.89	0	1,3,5	1.26	0
2	SAR	H	809	2	3,4,5	0.74	0	1,3,5	1.09	0
2	MVA	F	808	2	6,7,8	0.79	0	6,8,10	1.12	0
2	IML	H	811	2	7,8,9	0.83	0	6,9,11	1.17	1 (16%)
2	MVA	E	808	2	6,7,8	0.68	0	6,8,10	1.23	0
2	MVA	E	805	2	6,7,8	0.59	0	6,8,10	1.02	0
2	MVA	H	808	2	6,7,8	0.91	0	6,8,10	1.89	2 (33%)
2	MVA	G	808	2	6,7,8	0.72	0	6,8,10	0.98	0
2	SAR	E	809	2	3,4,5	0.83	0	1,3,5	0.88	0
2	MVA	E	807	2	6,7,8	0.97	0	6,8,10	2.32	2 (33%)
2	MVA	F	805	2	6,7,8	0.59	0	6,8,10	0.87	0
2	MVA	H	805	2	6,7,8	0.51	0	6,8,10	0.87	0
2	MVA	H	810	2	6,7,8	0.68	0	6,8,10	1.31	1 (16%)
2	MVA	G	805	2	6,7,8	0.47	0	6,8,10	0.97	0
2	MVA	H	807	2	6,7,8	1.25	1 (16%)	6,8,10	0.69	0

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	MVA	F	807	2	-	5/6/8/10	-
2	MVA	G	807	2	-	1/6/8/10	-
2	SAR	G	809	2	-	1/1/2/3	-
2	SAR	H	809	2	-	1/1/2/3	-
2	MVA	F	808	2	-	3/6/8/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IML	H	811	2	-	1/8/10/12	-
2	MVA	E	808	2	-	5/6/8/10	-
2	MVA	E	805	2	-	4/6/8/10	-
2	MVA	H	808	2	-	2/6/8/10	-
2	MVA	G	808	2	-	1/6/8/10	-
2	SAR	E	809	2	-	1/1/2/3	-
2	MVA	E	807	2	-	5/6/8/10	-
2	MVA	F	805	2	-	1/6/8/10	-
2	MVA	H	805	2	-	2/6/8/10	-
2	MVA	H	810	2	-	1/6/8/10	-
2	MVA	G	805	2	-	4/6/8/10	-
2	MVA	H	807	2	-	2/6/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	807	MVA	CA-N	2.63	1.51	1.47

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	807	MVA	CB-CA-C	-4.38	107.32	112.96
2	G	807	MVA	CB-CA-N	3.83	116.16	111.17
2	H	808	MVA	CB-CA-N	-3.80	106.22	111.17
2	G	807	MVA	O-C-CA	-2.85	117.32	124.86
2	G	807	MVA	CB-CA-C	-2.77	109.38	112.96
2	E	807	MVA	CG1-CB-CA	2.64	115.21	111.18
2	F	807	MVA	O-C-CA	-2.63	117.89	124.86
2	H	811	IML	CB-CA-C	-2.63	109.16	112.77
2	H	810	MVA	O-C-CA	-2.56	118.09	124.86
2	H	808	MVA	CB-CA-C	-2.14	110.20	112.96
2	F	807	MVA	CB-CA-C	-2.07	110.28	112.96

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	805	MVA	N-CA-CB-CG1
2	E	805	MVA	N-CA-CB-CG2
2	E	805	MVA	C-CA-CB-CG1

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Mol	Chain	Res	Type	Atoms
2	E	805	MVA	C-CA-CB-CG2
2	F	805	MVA	N-CA-CB-CG2
2	G	805	MVA	N-CA-CB-CG1
2	G	805	MVA	N-CA-CB-CG2
2	G	805	MVA	C-CA-CB-CG1
2	G	805	MVA	C-CA-CB-CG2
2	H	805	MVA	N-CA-CB-CG2
2	E	807	MVA	CB-CA-N-CN
2	E	807	MVA	N-CA-CB-CG1
2	E	807	MVA	N-CA-CB-CG2
2	E	807	MVA	C-CA-CB-CG1
2	E	807	MVA	C-CA-CB-CG2
2	F	807	MVA	CB-CA-N-CN
2	F	807	MVA	N-CA-CB-CG1
2	F	807	MVA	C-CA-CB-CG1
2	G	807	MVA	CB-CA-N-CN
2	H	807	MVA	CB-CA-N-CN
2	H	807	MVA	C-CA-CB-CG1
2	E	808	MVA	N-CA-CB-CG1
2	E	808	MVA	N-CA-CB-CG2
2	E	808	MVA	C-CA-CB-CG2
2	F	808	MVA	CB-CA-N-CN
2	F	808	MVA	N-CA-CB-CG1
2	F	808	MVA	N-CA-CB-CG2
2	G	808	MVA	CB-CA-N-CN
2	H	808	MVA	CB-CA-N-CN
2	E	809	SAR	C-CA-N-CN
2	G	809	SAR	C-CA-N-CN
2	H	809	SAR	C-CA-N-CN
2	H	810	MVA	CB-CA-N-CN
2	H	811	IML	CB-CA-N-CN
2	F	807	MVA	C-CA-CB-CG2
2	H	808	MVA	C-CA-CB-CG2
2	F	807	MVA	N-CA-CB-CG2
2	E	808	MVA	O-C-CA-CB
2	E	808	MVA	CB-CA-N-CN
2	H	805	MVA	N-CA-CB-CG1

There are no ring outliers.

16 monomers are involved in 44 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	807	MVA	4	0
2	G	807	MVA	2	0
2	G	809	SAR	4	0
2	H	809	SAR	4	0
2	F	808	MVA	1	0
2	H	811	IML	1	0
2	E	808	MVA	3	0
2	E	805	MVA	1	0
2	H	808	MVA	9	0
2	G	808	MVA	4	0
2	E	809	SAR	1	0
2	E	807	MVA	10	0
2	H	805	MVA	1	0
2	H	810	MVA	2	0
2	G	805	MVA	1	0
2	H	807	MVA	6	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 6 are monoatomic - leaving 9 for Mogul analysis.

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	$\# Z > 2$
5	PEG	D	807	-	6,6,6	0.18	0	5,5,5	0.12	0
3	EDO	C	806	-	3,3,3	0.09	0	2,2,2	0.17	0
5	PEG	C	802	-	6,6,6	0.24	0	5,5,5	0.11	0
3	EDO	D	802	-	3,3,3	0.06	0	2,2,2	0.18	0
3	EDO	C	801	-	3,3,3	0.11	0	2,2,2	0.21	0
3	EDO	D	801	-	3,3,3	0.09	0	2,2,2	0.08	0
4	BCT	B	802	-	3,3,3	1.34	0	2,3,3	1.22	0
3	EDO	B	801	-	3,3,3	0.05	0	2,2,2	0.04	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	D	803	-	3,3,3	0.13	0	2,2,2	0.23	0

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
5	PEG	D	807	-	-	2/4/4/4	-
3	EDO	C	806	-	-	0/1/1/1	-
5	PEG	C	802	-	-	1/4/4/4	-
3	EDO	D	802	-	-	1/1/1/1	-
3	EDO	C	801	-	-	1/1/1/1	-
3	EDO	D	801	-	-	1/1/1/1	-
3	EDO	B	801	-	-	1/1/1/1	-
3	EDO	D	803	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	807	PEG	C4-C3-O2-C2
5	D	807	PEG	O2-C3-C4-O4
3	D	802	EDO	O1-C1-C2-O2
3	D	801	EDO	O1-C1-C2-O2
3	C	801	EDO	O1-C1-C2-O2
3	D	803	EDO	O1-C1-C2-O2
5	C	802	PEG	C1-C2-O2-C3
3	B	801	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	720/745 (96%)	0.12	6 (0%) 82 80	54, 74, 98, 131	0
1	B	718/745 (96%)	0.29	6 (0%) 82 80	59, 79, 112, 135	0
1	C	706/745 (94%)	0.38	16 (2%) 61 58	52, 80, 117, 145	0
1	D	719/745 (96%)	0.09	7 (0%) 79 77	48, 68, 95, 112	0
2	E	1/15 (6%)	4.09	1 (100%) 0 0	112, 112, 112, 112	0
2	F	4/15 (26%)	2.07	2 (50%) 0 0	124, 125, 126, 131	0
2	G	1/15 (6%)	2.01	1 (100%) 0 0	107, 107, 107, 107	0
2	H	5/15 (33%)	1.17	1 (20%) 3 2	93, 93, 95, 98	0
All	All	2874/3040 (94%)	0.22	40 (1%) 73 71	48, 75, 111, 145	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	806	ILE	4.1
1	C	189	GLY	4.0
1	C	183	LEU	3.5
1	A	170	TRP	3.5
1	A	700	GLY	3.4
1	C	664	TYR	3.1
1	B	170	TRP	3.1
1	D	703	ALA	3.0
2	F	806	ILE	2.8
1	B	704	GLY	2.7
1	A	3	PHE	2.7
1	D	81	TRP	2.6
1	C	163	VAL	2.6
1	B	664	TYR	2.6
1	D	650	GLY	2.6
1	C	191	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	167	GLY	2.5
1	C	4	PRO	2.5
2	H	804	TRP	2.4
1	C	161	TYR	2.4
1	D	653	VAL	2.4
1	A	536	TYR	2.4
1	D	664	TYR	2.4
1	C	184	PRO	2.3
2	F	804	TRP	2.4
1	D	704	GLY	2.2
1	B	703	ALA	2.2
1	C	696	ASP	2.2
1	A	664	TYR	2.2
1	A	421	PHE	2.2
1	C	216	PHE	2.2
1	D	39	LEU	2.1
1	C	166	HIS	2.1
1	C	141	LEU	2.1
1	B	698	ASN	2.1
1	C	731	THR	2.1
1	B	728	THR	2.0
1	C	5	GLY	2.0
1	C	193	GLY	2.0
2	G	806	ILE	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MVA	G	805	8/9	0.54	0.24	101,104,108,108	0
2	MVA	E	805	8/9	0.58	0.22	105,109,110,110	0
2	MVA	H	810	8/9	0.80	0.15	100,101,102,103	0
2	MVA	E	808	8/9	0.84	0.17	104,106,109,109	0
2	MVA	G	808	8/9	0.84	0.16	108,109,110,112	0
2	MVA	F	805	8/9	0.84	0.20	122,124,126,127	0
2	SAR	E	809	5/6	0.85	0.17	98,102,105,105	0
2	MVA	F	808	8/9	0.85	0.24	106,113,116,117	0
2	SAR	G	809	5/6	0.86	0.22	92,97,103,104	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MVA	H	807	8/9	0.86	0.18	93,96,97,98	0
2	MVA	E	807	8/9	0.87	0.20	109,111,113,113	0
2	IML	H	811	9/10	0.87	0.17	92,94,98,98	0
2	SAR	H	809	5/6	0.89	0.12	99,99,100,102	0
2	MVA	F	807	8/9	0.89	0.22	121,122,124,125	0
2	MVA	G	807	8/9	0.89	0.20	106,107,108,109	0
2	MVA	H	808	8/9	0.90	0.14	97,99,99,100	0
2	MVA	H	805	8/9	0.94	0.15	93,95,98,99	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	D	801	4/4	0.77	0.13	77,79,81,83	0
3	EDO	D	803	4/4	0.77	0.13	77,77,77,78	0
6	NA	D	806	1/1	0.78	0.25	58,58,58,58	0
3	EDO	C	806	4/4	0.79	0.15	82,84,85,85	0
3	EDO	D	802	4/4	0.80	0.20	82,85,87,91	0
4	BCT	B	802	4/4	0.81	0.18	83,83,85,87	0
5	PEG	D	807	7/7	0.82	0.17	101,102,104,105	0
5	PEG	C	802	7/7	0.84	0.12	88,88,89,90	0
6	NA	D	804	1/1	0.85	0.13	62,62,62,62	0
3	EDO	B	801	4/4	0.85	0.15	81,81,83,83	0
3	EDO	C	801	4/4	0.86	0.15	74,74,75,75	0
6	NA	C	803	1/1	0.87	0.11	60,60,60,60	0
6	NA	D	805	1/1	0.89	0.14	64,64,64,64	0
6	NA	C	805	1/1	0.92	0.28	63,63,63,63	0
6	NA	C	804	1/1	0.92	0.13	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.