



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:16 PM GMT

PDB ID : 4L3T
Title : Crystal Structure of Substrate-free Human Presequence Protease
Authors : King, J.V.; Liang, W.G.; Tang, W.J.
Deposited on : 2013-06-06
Resolution : 2.03 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

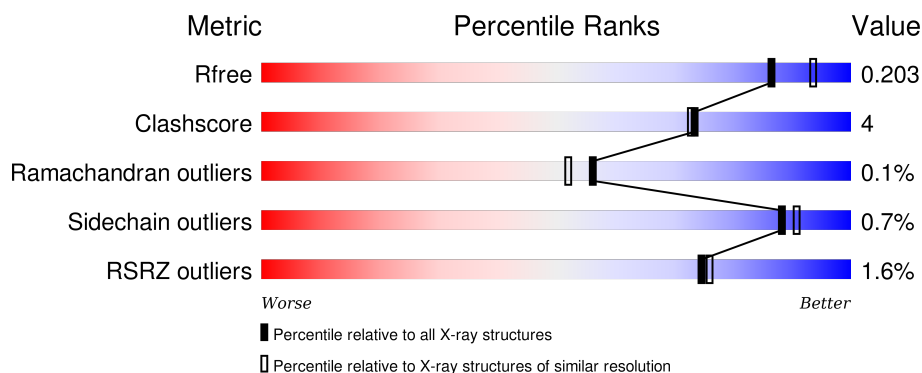
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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	7858 (2.04-2.00)
Clashscore	102246	9060 (2.04-2.00)
Ramachandran outliers	100387	8952 (2.04-2.00)
Sidechain outliers	100360	8951 (2.04-2.00)
RSRZ outliers	91569	7873 (2.04-2.00)

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	1014	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; width: 100%;"></div> <div style="position: absolute; bottom: -10px; left: 0; width: 100%;"></div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background-color: red;"></div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background-color: grey;"></div> </div> <div style="display: flex; justify-content: space-between; width: 100%;"> 89% 8% . </div> </div>
2	B	1014	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; width: 100%;"></div> <div style="position: absolute; bottom: -10px; left: 0; width: 100%;"></div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background-color: red;"></div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background-color: grey;"></div> </div> <div style="display: flex; justify-content: space-between; width: 100%;"> 87% 9% . </div> </div>

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
1	CAS	A	313	-	-	X	-
1	CAS	A	556	-	-	X	-
2	CAS	B	112	-	-	X	-
2	CAS	B	313	-	-	X	-
4	GOL	A	1107	-	-	-	X
4	GOL	B	1102	-	-	-	X
4	GOL	B	1112	-	-	-	X
5	ACT	A	1104	-	-	-	X
5	ACT	A	1105	-	-	-	X
5	ACT	A	1106	-	-	-	X
5	ACT	A	1108	-	-	-	X
5	ACT	B	1105	-	-	-	X
5	ACT	B	1106	-	-	-	X
5	ACT	B	1107	-	-	-	X
5	ACT	B	1109	-	-	-	X
5	ACT	B	1110	-	-	-	X
5	ACT	B	1113	-	-	-	X
5	ACT	B	1114	-	-	-	X

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 17046 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 Presequence protease, mitochondrial.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	As	C	N	O	S			
1	A	981	7977	6	5098	1349	1484	40	0	7	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	MET	-	EXPRESSION TAG	UNP Q5JRX3
A	25	HIS	-	EXPRESSION TAG	UNP Q5JRX3
A	26	HIS	-	EXPRESSION TAG	UNP Q5JRX3
A	27	HIS	-	EXPRESSION TAG	UNP Q5JRX3
A	28	HIS	-	EXPRESSION TAG	UNP Q5JRX3
A	29	HIS	-	EXPRESSION TAG	UNP Q5JRX3
A	30	HIS	-	EXPRESSION TAG	UNP Q5JRX3
A	31	ALA	-	EXPRESSION TAG	UNP Q5JRX3
A	32	ALA	-	EXPRESSION TAG	UNP Q5JRX3
A	107	GLN	GLU	ENGINEERED MUTATION	UNP Q5JRX3
A	328	VAL	ILE	SEE REMARK 999	UNP Q5JRX3
A	397	VAL	ALA	SEE REMARK 999	UNP Q5JRX3
A	1037	ARG	GLN	SEE REMARK 999	UNP Q5JRX3

- Molecule 2 is a protein called Presequence protease, mitochondrial.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	As	C	N	O	S			
2	B	978	7915	6	5064	1331	1474	40	0	3	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	24	MET	-	EXPRESSION TAG	UNP Q5JRX3
B	25	HIS	-	EXPRESSION TAG	UNP Q5JRX3
B	26	HIS	-	EXPRESSION TAG	UNP Q5JRX3

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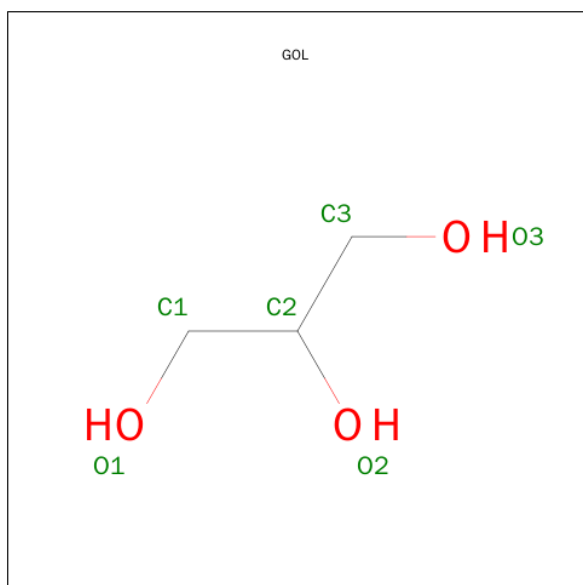
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Chain	Residue	Modelled	Actual	Comment	Reference
B	27	HIS	-	EXPRESSION TAG	UNP Q5JRX3
B	28	HIS	-	EXPRESSION TAG	UNP Q5JRX3
B	29	HIS	-	EXPRESSION TAG	UNP Q5JRX3
B	30	HIS	-	EXPRESSION TAG	UNP Q5JRX3
B	31	ALA	-	EXPRESSION TAG	UNP Q5JRX3
B	32	ALA	-	EXPRESSION TAG	UNP Q5JRX3
B	107	GLN	GLU	ENGINEERED MUTATION	UNP Q5JRX3
B	328	VAL	ILE	SEE REMARK 999	UNP Q5JRX3
B	397	VAL	ALA	SEE REMARK 999	UNP Q5JRX3
B	1037	ARG	GLN	SEE REMARK 999	UNP Q5JRX3

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Zn 1 1	0	0
3	A	1	Total Zn 1 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	553	Total 553	O 553	0	0
6	B	505	Total 505	O 505	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	245.78 Å 85.09 Å 158.46 Å 90.00° 127.54° 90.00°	Depositor
Resolution (Å)	42.55 – 2.03 42.55 – 2.03	Depositor EDS
% Data completeness (in resolution range)	89.3 (42.55-2.03) 84.7 (42.55-2.03)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.67 (at 2.03 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.171 , 0.210 0.168 , 0.203	Depositor DCC
R_{free} test set	7469 reflections (5.52%)	DCC
Wilson B-factor (Å ²)	27.2	Xtriage
Anisotropy	0.206	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.0	EDS
Estimated twinning fraction	0.013 for -h-2*k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 149408 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17046	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, CAS, ACT, MLZ, MLY

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	0.28	0/7704	0.47	1/10467 (0.0%)
2	B	0.28	0/7607	0.45	0/10329
All	All	0.28	0/15311	0.46	1/20796 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	316	ASP	CB-CG-OD2	5.17	122.95	118.30

There are no chirality outliers.

There are no planarity outliers.

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	7977	0	7860	57	1
2	B	7915	0	7797	74	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	18	0	24	2	0
4	B	24	0	32	1	1
5	A	16	0	12	0	0
5	B	36	0	27	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	553	0	0	8	4
6	B	505	0	0	7	4
All	All	17046	0	15752	131	5

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 (131) 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:A:118:PRO:HG2	1:A:556:CAS:CE2	1.51	1.40
2:B:609:GLU:CD	2:B:814:ARG:NH1	1.78	1.35
2:B:609:GLU:OE2	2:B:814:ARG:NH1	1.58	1.31
1:A:313:CAS:CE1	1:A:506:PRO:HB3	1.65	1.26
2:B:609:GLU:OE1	2:B:814:ARG:NH1	1.66	1.24
1:A:118:PRO:CG	1:A:556:CAS:CE2	2.30	1.09
2:B:609:GLU:CD	2:B:814:ARG:HH12	1.48	1.07
2:B:112:CAS:CE2	2:B:121:ASP:N	2.19	1.05
2:B:313:CAS:CE2	2:B:506:PRO:HG3	1.88	1.02
2:B:125:LYS:NZ	6:B:1556:HOH:O	1.93	1.00
2:B:112:CAS:CE2	2:B:121:ASP:H	1.74	0.98
2:B:313:CAS:SG	2:B:506:PRO:HA	2.06	0.95
2:B:112:CAS:CE2	2:B:121:ASP:CA	2.52	0.86
1:A:313:CAS:CE2	1:A:504:MET:HE2	2.05	0.85
1:A:313:CAS:CE1	1:A:506:PRO:CB	2.56	0.80
1:A:696:GLU:HG3	1:A:758:ILE:HD11	1.64	0.79
2:B:112:CAS:CE2	2:B:112:CAS:HA	2.14	0.77
1:A:313:CAS:CE2	1:A:504:MET:CE	2.63	0.77
1:A:913:SER:OG	1:A:915[A]:ASN:OD1	2.03	0.76
1:A:118:PRO:CD	1:A:556:CAS:CE2	2.64	0.75
2:B:171:CAS:HA	2:B:171:CAS:CE2	2.18	0.74
2:B:417:GLU:HG3	2:B:418:LYS:HG3	1.70	0.72
2:B:633:ARG:NH1	6:B:1567:HOH:O	2.23	0.72
2:B:112:CAS:CE2	2:B:121:ASP:HA	2.20	0.71
2:B:313:CAS:CE1	6:B:1689:HOH:O	2.40	0.70
1:A:121:ASP:OD1	6:A:1449:HOH:O	2.10	0.69
1:A:728:ARG:NH1	6:A:1282:HOH:O	2.26	0.68
1:A:666:VAL:HG21	1:A:772:LEU:HD21	1.76	0.67
2:B:119:CYS:HB2	6:B:1556:HOH:O	1.94	0.67
1:A:302:PRO:HA	1:A:498:HIS:HD2	1.58	0.67
2:B:324:MLZ:N	6:B:1597:HOH:O	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:823:PRO:HB3	1:A:982:SER:HB3	1.77	0.66
2:B:297:VAL:O	2:B:385:ARG:NH1	2.26	0.66
2:B:313:CAS:CE2	2:B:506:PRO:CG	2.71	0.65
2:B:690:ASN:ND2	5:B:1110:ACT:O	2.28	0.65
2:B:302:PRO:HA	2:B:498:HIS:HD2	1.62	0.64
1:A:845:GLU:OE2	6:A:1421:HOH:O	2.15	0.64
1:A:324:LYS:HG3	1:A:325:GLN:HG2	1.81	0.62
2:B:769:MLZ:HG2	2:B:773:LEU:HD12	1.81	0.62
1:A:118:PRO:HD2	1:A:556:CAS:CE2	2.30	0.61
2:B:313:CAS:SG	2:B:506:PRO:CA	2.84	0.60
2:B:604:LEU:HD13	2:B:772:LEU:HD21	1.83	0.60
2:B:580:ALA:N	2:B:581:GLY:HA2	2.18	0.59
2:B:75:ARG:NH2	2:B:458:ASP:OD1	2.36	0.58
2:B:303:TRP:H	2:B:498:HIS:CD2	2.21	0.58
2:B:728:ARG:NH2	2:B:744:ASP:OD2	2.35	0.58
1:A:791:GLN:OE1	6:A:1492:HOH:O	2.17	0.58
2:B:629:LEU:O	4:B:1102:GOL:O3	2.21	0.57
2:B:260:PRO:HB3	2:B:285:LEU:HD22	1.86	0.56
1:A:357:PRO:HA	1:A:362:TYR:CD1	2.41	0.56
1:A:677:LEU:HD21	1:A:792:THR:HA	1.87	0.56
2:B:313:CAS:HB2	2:B:504:MET:SD	2.46	0.55
2:B:313:CAS:AS	2:B:314:GLY:O	2.84	0.55
1:A:75:ARG:NH2	1:A:458:ASP:OD1	2.29	0.55
2:B:945:GLY:HA3	2:B:1002:LEU:HD21	1.89	0.55
1:A:128:ASN:OD1	1:A:888[B]:ARG:NH2	2.36	0.55
2:B:677:LEU:HD21	2:B:792:THR:HA	1.89	0.54
1:A:272:PRO:HA	4:A:1107:GOL:H2	1.88	0.54
2:B:109:THR:O	2:B:112:CAS:HB3	2.08	0.54
1:A:326:THR:HG21	1:A:401:ILE:HD11	1.89	0.54
2:B:130:SER:HB2	2:B:155:ASP:OD1	2.07	0.54
2:B:208:GLY:HA3	2:B:924:ARG:HD3	1.90	0.54
2:B:720:LEU:HD23	2:B:911:MLY:HH21	1.90	0.53
2:B:654:PRO:HB3	2:B:915[B]:ASN:HB3	1.91	0.53
1:A:313:CAS:CE2	1:A:504:MET:HE3	2.39	0.52
2:B:300:GLN:N	2:B:386:GLU:OE2	2.29	0.51
1:A:674:ASP:O	6:A:1492:HOH:O	2.19	0.51
2:B:112:CAS:CE2	2:B:112:CAS:CA	2.86	0.50
1:A:36:ARG:NH2	4:A:1107:GOL:O3	2.35	0.50
2:B:913:SER:OG	2:B:915[B]:ASN:OD1	2.15	0.49
1:A:389:PHE:CD1	1:A:493:PHE:HE2	2.30	0.49
2:B:398:GLU:OE2	2:B:505:ARG:NE	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:603:SER:HB3	1:A:777[A]:ASN:HB3	1.95	0.49
1:A:119:CYS:N	1:A:556:CAS:CE2	2.76	0.49
1:A:396:ILE:HD11	1:A:401:ILE:HG12	1.95	0.48
2:B:389:PHE:CD2	2:B:493:PHE:HE2	2.32	0.47
1:A:699:PHE:HD2	1:A:758:ILE:HD13	1.79	0.47
2:B:960:PHE:HA	2:B:963:VAL:HG22	1.97	0.47
2:B:303:TRP:H	2:B:498:HIS:HD2	1.61	0.47
2:B:814:ARG:C	2:B:816:HIS:N	2.68	0.46
2:B:654:PRO:CB	2:B:915[B]:ASN:HB3	2.46	0.46
2:B:357:PRO:HA	2:B:362:TYR:CD1	2.50	0.46
1:A:325:GLN:HG3	1:A:506:PRO:HG3	1.98	0.46
1:A:363:MLY:HA	1:A:367:GLU:HG3	1.98	0.45
2:B:844:MET:HG2	2:B:846:PRO:HD3	1.97	0.45
1:A:364:ALA:O	1:A:368:SER:HB3	2.16	0.45
1:A:363:MLY:O	1:A:367:GLU:HB2	2.16	0.45
2:B:776:ASP:OD1	2:B:806:SER:N	2.48	0.45
2:B:397:VAL:HG12	2:B:399:LYS:HG2	1.99	0.45
2:B:690:ASN:ND2	6:B:1564:HOH:O	2.49	0.45
2:B:612:ARG:NH2	2:B:664:GLN:OE1	2.47	0.45
1:A:625:LEU:O	1:A:692:CAS:HB3	2.18	0.44
1:A:155:ASP:HB2	1:A:560:LEU:HD11	2.00	0.44
2:B:695:GLU:HG3	2:B:698:HIS:HB3	1.98	0.43
1:A:750:MLY:HD3	1:A:750:MLY:HH12	1.75	0.43
2:B:661:THR:HG22	2:B:842:LEU:HD23	2.00	0.43
1:A:313:CAS:SG	1:A:506:PRO:HA	2.58	0.43
2:B:914:HIS:O	6:B:1425:HOH:O	2.21	0.43
1:A:100:THR:HB	1:A:245:LEU:HB2	2.01	0.43
2:B:578:LEU:HA	2:B:578:LEU:HD13	1.88	0.43
1:A:599:ARG:HA	1:A:668:PHE:O	2.19	0.42
2:B:609:GLU:OE2	2:B:814:ARG:HD2	2.19	0.42
1:A:289:GLN:NE2	6:A:1442:HOH:O	2.32	0.42
2:B:816:HIS:O	2:B:843:VAL:HA	2.19	0.42
1:A:543:GLU:O	1:A:547[B]:GLN:HG2	2.20	0.42
1:A:38:LEU:HD21	1:A:56:SER:OG	2.20	0.42
2:B:302:PRO:HA	2:B:498:HIS:CD2	2.50	0.42
1:A:367:GLU:OE2	6:A:1587:HOH:O	2.20	0.42
1:A:174:GLU:OE1	1:A:540:MLY:HH23	2.20	0.42
1:A:654:PRO:CG	1:A:915[A]:ASN:HB3	2.50	0.42
2:B:888:ARG:HD2	2:B:888:ARG:HA	1.79	0.42
2:B:278:MLY:HH13	2:B:278:MLY:HD2	1.77	0.42
2:B:858:LEU:HD13	2:B:1036:ILE:HG23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:305:LYS:HA	2:B:306:PRO:HD3	1.95	0.41
2:B:549:SER:OG	2:B:902:MLY:HH13	2.19	0.41
2:B:1000:MLZ:HD3	2:B:1000:MLZ:HCM3	1.85	0.41
1:A:105:ILE:HG21	1:A:177:PHE:HE1	1.85	0.41
1:A:66:MLY:HH12	1:A:460:ASP:OD1	2.20	0.41
1:A:769:MLZ:HG2	1:A:773:LEU:HD12	2.03	0.41
1:A:700:MLY:CG	1:A:758:ILE:HG12	2.50	0.41
1:A:123:PHE:HB3	6:A:1449:HOH:O	2.19	0.41
2:B:31:ALA:HA	2:B:32:ALA:HA	1.79	0.41
2:B:772:LEU:HD23	2:B:772:LEU:HA	1.91	0.41
1:A:208:GLY:HA3	1:A:924:ARG:HD3	2.02	0.41
2:B:466:MLY:HH22	2:B:466:MLY:HD3	1.85	0.41
2:B:154:LYS:HD2	2:B:154:LYS:HA	1.86	0.40
2:B:174:GLU:OE1	2:B:540:MLY:HE3	2.22	0.40
2:B:301:THR:HA	2:B:302:PRO:HD3	1.96	0.40
1:A:777[A]:ASN:OD1	1:A:841:LYS:HD2	2.21	0.40
1:A:115:GLN:NE2	1:A:171:CAS:CE1	2.84	0.40
1:A:554:ALA:C	1:A:556:CAS:N	2.75	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:1593:HOH:O	6:B:1534:HOH:O[3_556]	1.96	0.24
6:A:1557:HOH:O	6:B:1550:HOH:O[4_554]	2.10	0.10
1:A:682:GLN:NE2	4:B:1102:GOL:O3[3_556]	2.13	0.07
6:A:1601:HOH:O	6:B:1564:HOH:O[3_556]	2.16	0.04
6:A:1557:HOH:O	6:B:1548:HOH:O[4_554]	2.19	0.01

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	937/1014 (92%)	919 (98%)	17 (2%)	1 (0%)	56	52
2	B	926/1014 (91%)	909 (98%)	16 (2%)	1 (0%)	56	52
All	All	1863/2028 (92%)	1828 (98%)	33 (2%)	2 (0%)	56	52

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	147	PRO
2	B	147	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	838/858 (98%)	833 (99%)	5 (1%)	90	92
2	B	827/855 (97%)	819 (99%)	8 (1%)	82	85
All	All	1665/1713 (97%)	1652 (99%)	13 (1%)	88	89

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	292	GLU
1	A	728	ARG
1	A	777[A]	ASN
1	A	777[B]	ASN
1	A	981	LEU
2	B	75	ARG
2	B	294	SER
2	B	403	THR
2	B	604	LEU
2	B	728	ARG
2	B	814	ARG
2	B	843	VAL
2	B	1011	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	498	HIS
1	A	989	HIS
2	B	498	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

89 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	MLY	A	1000	1	9,10,11	0.52	0	9,11,13	0.84	0
1	MLY	A	1013	1	9,10,11	0.51	0	9,11,13	0.81	0
1	MLY	A	116	1	9,10,11	0.44	0	9,11,13	1.17	1 (11%)
1	MLY	A	154	1	9,10,11	0.53	0	9,11,13	0.95	0
1	CAS	A	171	1	5,8,9	1.07	0	2,9,11	2.04	1 (50%)
1	MLZ	A	207	1	8,9,10	1.07	1 (12%)	7,9,11	1.20	0
1	CAS	A	241	1	5,8,9	1.09	0	2,9,11	1.92	1 (50%)
1	MLY	A	251	1	9,10,11	0.55	0	9,11,13	0.94	0
1	MLY	A	278	1	9,10,11	0.47	0	9,11,13	0.88	0
1	MLY	A	287	1	9,10,11	0.45	0	9,11,13	0.96	0
1	MLY	A	290	1	9,10,11	0.51	0	9,11,13	0.86	0
1	CAS	A	313	1	5,8,9	1.07	0	2,9,11	1.68	1 (50%)
1	MLY	A	363	1	9,10,11	0.54	0	9,11,13	0.86	0
1	MLY	A	431	1	9,10,11	0.55	0	9,11,13	0.90	0
1	MLZ	A	437	1	8,9,10	0.91	0	7,9,11	1.20	1 (14%)
1	MLY	A	466	1	9,10,11	0.49	0	9,11,13	0.96	0
1	CAS	A	477	1	5,8,9	1.08	0	2,9,11	1.70	1 (50%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	MLY	A	488	1	9,10,11	0.52	0	9,11,13	0.78	0
1	MLY	A	513	1	9,10,11	0.48	0	9,11,13	0.83	0
1	MLY	A	521	1	9,10,11	0.50	0	9,11,13	0.86	0
1	MLZ	A	525	1	8,9,10	0.96	0	7,9,11	1.14	0
1	MLY	A	540	1	9,10,11	0.50	0	9,11,13	0.83	0
1	MLZ	A	550	1	8,9,10	0.92	0	7,9,11	1.24	1 (14%)
1	CAS	A	556	1	5,8,9	1.09	0	2,9,11	1.89	1 (50%)
1	MLY	A	642	1	9,10,11	0.49	0	9,11,13	1.00	1 (11%)
1	MLY	A	66	1	9,10,11	0.51	0	9,11,13	0.90	0
1	CAS	A	692	1	5,8,9	1.06	0	2,9,11	1.81	0
1	MLY	A	700	1	9,10,11	0.51	0	9,11,13	0.88	0
1	MLY	A	704	1	9,10,11	0.51	0	9,11,13	0.94	1 (11%)
1	MLY	A	750	1	9,10,11	0.51	0	9,11,13	0.83	0
1	MLY	A	759	1	9,10,11	0.50	0	9,11,13	0.99	1 (11%)
1	MLY	A	764	1	9,10,11	0.50	0	9,11,13	0.93	0
1	MLZ	A	769	1	8,9,10	0.77	0	7,9,11	0.91	0
1	MLY	A	794	1	9,10,11	0.58	0	9,11,13	0.80	0
1	MLY	A	854	1	9,10,11	0.61	0	9,11,13	0.87	0
1	MLY	A	884	1	9,10,11	0.51	0	9,11,13	0.88	0
1	MLY	A	902	1	9,10,11	0.54	0	9,11,13	0.85	0
1	MLY	A	911	1	9,10,11	0.58	0	9,11,13	0.85	0
1	MLY	A	937	1	9,10,11	0.60	0	9,11,13	0.86	0
1	MLY	A	943	1	9,10,11	0.64	0	9,11,13	0.85	0
1	MLY	A	946	1	9,10,11	0.49	0	9,11,13	0.93	0
1	MLY	A	956	1	9,10,11	0.46	0	9,11,13	1.40	2 (22%)
1	MLY	A	972	1	9,10,11	0.54	0	9,11,13	0.96	1 (11%)
2	MLZ	B	1000	2	8,9,10	1.12	1 (12%)	7,9,11	1.11	0
2	MLY	B	1013	2	9,10,11	0.54	0	9,11,13	0.88	0
2	CAS	B	112	2	5,8,9	1.06	0	2,9,11	2.05	1 (50%)
2	MLY	B	116	2	9,10,11	0.55	0	9,11,13	0.91	0
2	CAS	B	171	2	5,8,9	1.07	0	2,9,11	2.06	1 (50%)
2	MLY	B	199	2	9,10,11	0.45	0	9,11,13	1.06	1 (11%)
2	MLY	B	251	2	9,10,11	0.48	0	9,11,13	0.92	0
2	MLY	B	278	2	9,10,11	0.49	0	9,11,13	0.96	0
2	MLY	B	287	2	9,10,11	0.50	0	9,11,13	1.00	1 (11%)
2	MLY	B	290	2	9,10,11	0.48	0	9,11,13	0.94	0
2	CAS	B	313	2	5,8,9	1.09	0	2,9,11	1.62	0
2	MLZ	B	324	2	8,9,10	0.97	0	7,9,11	1.08	0
2	MLY	B	363	2	9,10,11	0.64	0	9,11,13	0.79	0
2	MLY	B	41	2	9,10,11	0.53	0	9,11,13	1.02	1 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MLY	B	431	2	9,10,11	0.52	0	9,11,13	0.90	0
2	MLY	B	466	2	9,10,11	0.53	0	9,11,13	0.91	0
2	MLY	B	473	2	9,10,11	0.47	0	9,11,13	0.93	0
2	CAS	B	477	2	5,8,9	1.09	0	2,9,11	2.06	1 (50%)
2	MLY	B	488	2	9,10,11	0.53	0	9,11,13	0.87	0
2	MLZ	B	490	2	8,9,10	1.06	1 (12%)	7,9,11	1.19	0
2	MLZ	B	494	2	8,9,10	0.82	0	7,9,11	1.09	0
2	MLY	B	499	2	9,10,11	0.51	0	9,11,13	0.96	0
2	MLY	B	513	2	9,10,11	0.52	0	9,11,13	0.84	0
2	MLY	B	525	2	9,10,11	0.52	0	9,11,13	0.87	0
2	MLY	B	540	2	9,10,11	0.52	0	9,11,13	0.80	0
2	MLY	B	550	2	9,10,11	0.57	0	9,11,13	0.95	0
2	CAS	B	556	2	5,8,9	1.09	0	2,9,11	1.93	1 (50%)
2	MLY	B	624	2	9,10,11	0.44	0	9,11,13	1.04	1 (11%)
2	MLY	B	642	2	9,10,11	0.56	0	9,11,13	0.96	1 (11%)
2	MLY	B	66	2	9,10,11	0.55	0	9,11,13	0.95	0
2	CAS	B	692	2	5,8,9	1.07	0	2,9,11	1.81	1 (50%)
2	MLY	B	704	2	9,10,11	0.47	0	9,11,13	0.96	0
2	MLY	B	750	2	9,10,11	0.53	0	9,11,13	0.92	1 (11%)
2	MLZ	B	759	2	8,9,10	0.80	0	7,9,11	1.32	1 (14%)
2	MLZ	B	764	2	8,9,10	0.99	0	7,9,11	1.13	0
2	MLZ	B	769	2	8,9,10	0.88	0	7,9,11	1.13	1 (14%)
2	MLZ	B	854	2	8,9,10	0.96	0	7,9,11	1.12	1 (14%)
2	MLZ	B	884	2	8,9,10	0.91	0	7,9,11	1.06	0
2	MLY	B	902	2	9,10,11	0.54	0	9,11,13	0.88	0
2	MLY	B	911	2	9,10,11	0.47	0	9,11,13	0.97	0
2	MLZ	B	937	2	8,9,10	0.97	1 (12%)	7,9,11	1.18	1 (14%)
2	MLY	B	943	2	9,10,11	0.58	0	9,11,13	0.85	0
2	MLY	B	946	2	9,10,11	0.54	0	9,11,13	0.96	0
2	MLY	B	956	2	9,10,11	0.60	0	9,11,13	1.00	0
2	MLY	B	972	2	9,10,11	0.47	0	9,11,13	0.92	0
2	MLY	B	986	2	9,10,11	0.54	0	9,11,13	0.84	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
1	MLY	A	1000	1	-	0/7/9/11	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MLY	A	1013	1	-	0/7/9/11	0/0/0/0
1	MLY	A	116	1	-	0/7/9/11	0/0/0/0
1	MLY	A	154	1	-	0/7/9/11	0/0/0/0
1	CAS	A	171	1	-	0/0/7/9	0/0/0/0
1	MLZ	A	207	1	-	0/6/8/10	0/0/0/0
1	CAS	A	241	1	-	0/0/7/9	0/0/0/0
1	MLY	A	251	1	-	0/7/9/11	0/0/0/0
1	MLY	A	278	1	-	0/7/9/11	0/0/0/0
1	MLY	A	287	1	-	0/7/9/11	0/0/0/0
1	MLY	A	290	1	-	0/7/9/11	0/0/0/0
1	CAS	A	313	1	-	0/0/7/9	0/0/0/0
1	MLY	A	363	1	-	0/7/9/11	0/0/0/0
1	MLY	A	431	1	-	0/7/9/11	0/0/0/0
1	MLZ	A	437	1	-	0/6/8/10	0/0/0/0
1	MLY	A	466	1	-	0/7/9/11	0/0/0/0
1	CAS	A	477	1	-	0/0/7/9	0/0/0/0
1	MLY	A	488	1	-	0/7/9/11	0/0/0/0
1	MLY	A	513	1	-	0/7/9/11	0/0/0/0
1	MLY	A	521	1	-	0/7/9/11	0/0/0/0
1	MLZ	A	525	1	-	0/6/8/10	0/0/0/0
1	MLY	A	540	1	-	0/7/9/11	0/0/0/0
1	MLZ	A	550	1	-	0/6/8/10	0/0/0/0
1	CAS	A	556	1	-	0/0/7/9	0/0/0/0
1	MLY	A	642	1	-	0/7/9/11	0/0/0/0
1	MLY	A	66	1	-	0/7/9/11	0/0/0/0
1	CAS	A	692	1	-	0/0/7/9	0/0/0/0
1	MLY	A	700	1	-	0/7/9/11	0/0/0/0
1	MLY	A	704	1	-	0/7/9/11	0/0/0/0
1	MLY	A	750	1	-	0/7/9/11	0/0/0/0
1	MLY	A	759	1	-	0/7/9/11	0/0/0/0
1	MLY	A	764	1	-	0/7/9/11	0/0/0/0
1	MLZ	A	769	1	-	0/6/8/10	0/0/0/0
1	MLY	A	794	1	-	0/7/9/11	0/0/0/0
1	MLY	A	854	1	-	0/7/9/11	0/0/0/0
1	MLY	A	884	1	-	0/7/9/11	0/0/0/0
1	MLY	A	902	1	-	0/7/9/11	0/0/0/0
1	MLY	A	911	1	-	0/7/9/11	0/0/0/0
1	MLY	A	937	1	-	0/7/9/11	0/0/0/0
1	MLY	A	943	1	-	0/7/9/11	0/0/0/0
1	MLY	A	946	1	-	0/7/9/11	0/0/0/0
1	MLY	A	956	1	-	0/7/9/11	0/0/0/0
1	MLY	A	972	1	-	0/7/9/11	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MLZ	B	1000	2	-	0/6/8/10	0/0/0/0
2	MLY	B	1013	2	-	0/7/9/11	0/0/0/0
2	CAS	B	112	2	-	0/0/7/9	0/0/0/0
2	MLY	B	116	2	-	0/7/9/11	0/0/0/0
2	CAS	B	171	2	-	0/0/7/9	0/0/0/0
2	MLY	B	199	2	-	0/7/9/11	0/0/0/0
2	MLY	B	251	2	-	0/7/9/11	0/0/0/0
2	MLY	B	278	2	-	0/7/9/11	0/0/0/0
2	MLY	B	287	2	-	0/7/9/11	0/0/0/0
2	MLY	B	290	2	-	0/7/9/11	0/0/0/0
2	CAS	B	313	2	-	0/0/7/9	0/0/0/0
2	MLZ	B	324	2	-	0/6/8/10	0/0/0/0
2	MLY	B	363	2	-	0/7/9/11	0/0/0/0
2	MLY	B	41	2	-	0/7/9/11	0/0/0/0
2	MLY	B	431	2	-	0/7/9/11	0/0/0/0
2	MLY	B	466	2	-	0/7/9/11	0/0/0/0
2	MLY	B	473	2	-	0/7/9/11	0/0/0/0
2	CAS	B	477	2	-	0/0/7/9	0/0/0/0
2	MLY	B	488	2	-	0/7/9/11	0/0/0/0
2	MLZ	B	490	2	-	0/6/8/10	0/0/0/0
2	MLZ	B	494	2	-	0/6/8/10	0/0/0/0
2	MLY	B	499	2	-	0/7/9/11	0/0/0/0
2	MLY	B	513	2	-	0/7/9/11	0/0/0/0
2	MLY	B	525	2	-	0/7/9/11	0/0/0/0
2	MLY	B	540	2	-	0/7/9/11	0/0/0/0
2	MLY	B	550	2	-	0/7/9/11	0/0/0/0
2	CAS	B	556	2	-	0/0/7/9	0/0/0/0
2	MLY	B	624	2	-	0/7/9/11	0/0/0/0
2	MLY	B	642	2	-	0/7/9/11	0/0/0/0
2	MLY	B	66	2	-	0/7/9/11	0/0/0/0
2	CAS	B	692	2	-	0/0/7/9	0/0/0/0
2	MLY	B	704	2	-	0/7/9/11	0/0/0/0
2	MLY	B	750	2	-	0/7/9/11	0/0/0/0
2	MLZ	B	759	2	-	0/6/8/10	0/0/0/0
2	MLZ	B	764	2	-	0/6/8/10	0/0/0/0
2	MLZ	B	769	2	-	0/6/8/10	0/0/0/0
2	MLZ	B	854	2	-	0/6/8/10	0/0/0/0
2	MLZ	B	884	2	-	0/6/8/10	0/0/0/0
2	MLY	B	902	2	-	0/7/9/11	0/0/0/0
2	MLY	B	911	2	-	0/7/9/11	0/0/0/0
2	MLZ	B	937	2	-	0/6/8/10	0/0/0/0
2	MLY	B	943	2	-	0/7/9/11	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MLY	B	946	2	-	0/7/9/11	0/0/0/0
2	MLY	B	956	2	-	0/7/9/11	0/0/0/0
2	MLY	B	972	2	-	0/7/9/11	0/0/0/0
2	MLY	B	986	2	-	0/7/9/11	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1000	MLZ	CB-CA	-2.57	1.51	1.53
1	A	207	MLZ	CB-CA	-2.47	1.51	1.53
2	B	490	MLZ	CB-CA	-2.23	1.51	1.53
2	B	937	MLZ	CB-CA	-2.06	1.51	1.53

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	759	MLZ	O-C-CA	-2.71	118.44	125.49
2	B	171	CAS	O-C-CA	-2.59	118.73	125.49
2	B	477	CAS	O-C-CA	-2.58	118.78	125.49
1	A	171	CAS	O-C-CA	-2.55	118.84	125.49
1	A	956	MLY	CD-CE-NZ	-2.55	107.36	113.92
2	B	112	CAS	O-C-CA	-2.51	118.95	125.49
2	B	556	CAS	O-C-CA	-2.41	119.22	125.49
1	A	477	CAS	O-C-CA	-2.40	119.23	125.49
1	A	241	CAS	O-C-CA	-2.40	119.23	125.49
1	A	556	CAS	O-C-CA	-2.33	119.42	125.49
1	A	759	MLY	O-C-CA	-2.30	119.49	125.49
1	A	956	MLY	CH2-NZ-CH1	-2.29	103.60	109.72
1	A	550	MLZ	O-C-CA	-2.28	119.54	125.49
1	A	437	MLZ	O-C-CA	-2.21	119.73	125.49
2	B	854	MLZ	O-C-CA	-2.16	119.87	125.49
2	B	692	CAS	O-C-CA	-2.14	119.91	125.49
1	A	704	MLY	O-C-CA	-2.13	119.94	125.49
1	A	642	MLY	O-C-CA	-2.12	119.95	125.49
2	B	41	MLY	O-C-CA	-2.12	119.97	125.49
2	B	642	MLY	O-C-CA	-2.09	120.06	125.49
2	B	199	MLY	O-C-CA	-2.08	120.06	125.49
2	B	769	MLZ	O-C-CA	-2.07	120.10	125.49
2	B	624	MLY	O-C-CA	-2.03	120.20	125.49
2	B	937	MLZ	O-C-CA	-2.03	120.21	125.49
1	A	116	MLY	CD-CE-NZ	-2.02	108.71	113.92
2	B	750	MLY	O-C-CA	-2.02	120.22	125.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	313	CAS	O-C-CA	-2.01	120.25	125.49
1	A	972	MLY	O-C-CA	-2.01	120.25	125.49
2	B	287	MLY	O-C-CA	-2.01	120.26	125.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

21 monomers are involved in 44 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	171	CAS	1	0
1	A	313	CAS	6	0
1	A	363	MLY	2	0
1	A	540	MLY	1	0
1	A	556	CAS	6	0
1	A	66	MLY	1	0
1	A	692	CAS	1	0
1	A	700	MLY	1	0
1	A	750	MLY	1	0
1	A	769	MLZ	1	0
2	B	1000	MLZ	1	0
2	B	112	CAS	7	0
2	B	171	CAS	1	0
2	B	278	MLY	1	0
2	B	313	CAS	7	0
2	B	324	MLZ	1	0
2	B	466	MLY	1	0
2	B	540	MLY	1	0
2	B	769	MLZ	1	0
2	B	902	MLY	1	0
2	B	911	MLY	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 22 ligands modelled in this entry, 2 are monoatomic - leaving 20 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
4	GOL	A	1102	-	5,5,5	0.31	0	5,5,5	0.31	0
4	GOL	A	1103	-	5,5,5	0.28	0	5,5,5	0.32	0
5	ACT	A	1104	-	1,3,3	1.02	0	0,3,3	0.00	-
5	ACT	A	1105	-	1,3,3	1.13	0	0,3,3	0.00	-
5	ACT	A	1106	-	1,3,3	1.33	0	0,3,3	0.00	-
4	GOL	A	1107	-	5,5,5	0.32	0	5,5,5	0.42	0
5	ACT	A	1108	3	1,3,3	1.36	0	0,3,3	0.00	-
4	GOL	B	1102	-	5,5,5	0.36	0	5,5,5	0.50	0
4	GOL	B	1103	-	5,5,5	0.29	0	5,5,5	0.35	0
4	GOL	B	1104	-	5,5,5	0.34	0	5,5,5	0.10	0
5	ACT	B	1105	-	1,3,3	1.36	0	0,3,3	0.00	-
5	ACT	B	1106	-	1,3,3	1.14	0	0,3,3	0.00	-
5	ACT	B	1107	-	1,3,3	0.97	0	0,3,3	0.00	-
5	ACT	B	1108	-	1,3,3	1.21	0	0,3,3	0.00	-
5	ACT	B	1109	-	1,3,3	1.10	0	0,3,3	0.00	-
5	ACT	B	1110	-	1,3,3	1.18	0	0,3,3	0.00	-
5	ACT	B	1111	-	1,3,3	1.19	0	0,3,3	0.00	-
4	GOL	B	1112	-	5,5,5	0.33	0	5,5,5	0.29	0
5	ACT	B	1113	-	1,3,3	1.32	0	0,3,3	0.00	-
5	ACT	B	1114	-	1,3,3	0.90	0	0,3,3	0.00	-

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
4	GOL	A	1102	-	-	0/4/4/4	0/0/0/0
4	GOL	A	1103	-	-	0/4/4/4	0/0/0/0
5	ACT	A	1104	-	-	0/0/0/0	0/0/0/0
5	ACT	A	1105	-	-	0/0/0/0	0/0/0/0
5	ACT	A	1106	-	-	0/0/0/0	0/0/0/0
4	GOL	A	1107	-	-	0/4/4/4	0/0/0/0
5	ACT	A	1108	3	-	0/0/0/0	0/0/0/0
4	GOL	B	1102	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	1103	-	-	0/4/4/4	0/0/0/0
4	GOL	B	1104	-	-	0/4/4/4	0/0/0/0
5	ACT	B	1105	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1106	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1107	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1108	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1109	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1110	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1111	-	-	0/0/0/0	0/0/0/0
4	GOL	B	1112	-	-	0/4/4/4	0/0/0/0
5	ACT	B	1113	-	-	0/0/0/0	0/0/0/0
5	ACT	B	1114	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1107	GOL	2	0
4	B	1102	GOL	1	1
5	B	1110	ACT	1	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	938/1014 (92%)	-0.35	13 (1%) 78 79	18, 30, 58, 93	0
2	B	932/1014 (91%)	-0.42	16 (1%) 73 74	20, 32, 58, 98	0
All	All	1870/2028 (92%)	-0.39	29 (1%) 74 76	18, 31, 58, 98	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	838	VAL	6.3
2	B	838	VAL	5.8
1	A	316	ASP	4.9
1	A	811	ARG	4.6
2	B	31	ALA	3.9
2	B	816	HIS	3.8
1	A	531	PRO	3.7
1	A	823	PRO	3.5
2	B	314	GLY	3.3
1	A	571	VAL	3.0
2	B	138[A]	PHE	2.8
1	A	822	VAL	2.8
2	B	819	GLU	2.6
1	A	813	VAL	2.6
1	A	821	PRO	2.5
2	B	1030	LYS	2.5
2	B	813	VAL	2.5
1	A	508	ASP	2.4
1	A	529	LEU	2.4
2	B	399	LYS	2.4
2	B	983	ASP	2.3
2	B	812	PRO	2.3
1	A	527	GLU	2.3
2	B	817	THR	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	1027	LYS	2.3
2	B	851	TRP	2.2
2	B	839	ILE	2.1
2	B	756	THR	2.1
1	A	839	ILE	2.1

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. 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
2	MLY	B	525	11/12	0.95	0.12	-	25,39,47,51	0
2	CAS	B	556	9/10	0.81	0.18	-	34,42,114,157	0
2	MLY	B	750	11/12	0.93	0.14	-	24,36,65,65	0
1	MLY	A	540	11/12	0.95	0.10	-	28,30,45,47	0
1	MLY	A	956	11/12	0.96	0.12	-	18,22,62,69	0
1	MLY	A	488	11/12	0.96	0.09	-	28,35,39,40	0
1	CAS	A	313	9/10	0.67	0.25	-	54,56,93,287	0
1	MLY	A	764	11/12	0.97	0.11	-	35,38,49,50	0
1	CAS	A	241	9/10	0.94	0.13	-	22,39,71,190	0
1	CAS	A	556	9/10	0.76	0.23	-	24,32,90,102	0
2	MLY	B	473	11/12	0.96	0.12	-	31,37,68,69	0
1	MLY	A	884	11/12	0.98	0.09	-	20,25,46,51	0
1	MLZ	A	525	10/11	0.93	0.15	-	41,45,51,53	0
2	MLY	B	972	11/12	0.93	0.13	-	27,36,71,71	0
2	MLZ	B	759	10/11	0.96	0.13	-	41,47,64,65	0
2	MLY	B	466	11/12	0.96	0.10	-	26,36,47,49	0
1	MLY	A	902	11/12	0.97	0.12	-	25,27,42,43	0
1	MLZ	A	207	10/11	0.96	0.12	-	19,28,53,56	0
1	CAS	A	692	9/10	0.73	0.20	-	39,44,108,242	0
2	MLY	B	116	11/12	0.96	0.12	-	21,35,57,58	0
2	MLY	B	540	11/12	0.97	0.10	-	20,26,41,45	0
2	CAS	B	313	9/10	0.66	0.25	-	48,54,100,262	0
2	MLY	B	251	11/12	0.97	0.10	-	23,26,55,56	0
1	MLY	A	513	11/12	0.92	0.18	-	43,48,53,60	0
2	MLY	B	363	11/12	0.97	0.10	-	28,34,59,60	0
2	CAS	B	112	9/10	0.92	0.17	-	25,33,104,226	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CAS	B	477	9/10	0.94	0.11	-	35,47,118,187	0
1	MLY	A	759	11/12	0.95	0.17	-	32,42,63,64	0
2	MLY	B	1013	11/12	0.93	0.09	-	31,39,54,55	0
1	MLY	A	700	11/12	0.94	0.18	-	28,40,73,74	0
2	MLY	B	902	11/12	0.97	0.09	-	27,31,40,41	0
2	MLZ	B	854	10/11	0.96	0.07	-	29,33,46,48	0
2	MLZ	B	494	10/11	0.95	0.19	-	29,42,61,62	0
2	CAS	B	692	9/10	0.87	0.14	-	28,41,70,165	0
2	MLY	B	550	11/12	0.93	0.14	-	29,37,51,58	0
2	MLY	B	290	11/12	0.94	0.12	-	36,39,65,65	0
1	MLY	A	251	11/12	0.96	0.10	-	24,27,53,54	0
1	MLY	A	431	11/12	0.96	0.10	-	24,33,63,63	0
1	MLY	A	1013	11/12	0.93	0.16	-	35,38,56,58	0
1	MLY	A	521	11/12	0.92	0.22	-	44,59,87,89	0
2	MLZ	B	324	10/11	0.90	0.20	-	47,62,75,75	0
1	MLY	A	116	11/12	0.96	0.10	-	25,30,42,46	0
2	MLZ	B	490	10/11	0.97	0.14	-	31,37,66,68	0
2	MLY	B	513	11/12	0.86	0.15	-	46,49,53,56	0
2	MLY	B	41	11/12	0.97	0.10	-	27,29,44,49	0
2	MLZ	B	1000	10/11	0.97	0.12	-	25,31,64,65	0
1	MLY	A	290	11/12	0.96	0.11	-	34,39,59,60	0
1	MLY	A	972	11/12	0.93	0.14	-	22,35,63,64	0
2	MLY	B	278	11/12	0.97	0.10	-	24,29,60,62	0
1	MLY	A	1000	11/12	0.97	0.11	-	20,27,60,64	0
2	MLZ	B	764	10/11	0.97	0.09	-	35,44,48,50	0
2	MLY	B	66	11/12	0.97	0.13	-	20,24,58,59	0
1	MLY	A	943	11/12	0.95	0.11	-	20,29,43,47	0
1	CAS	A	171	9/10	0.88	0.15	-	19,34,86,100	0
1	MLZ	A	769	10/11	0.98	0.08	-	27,31,43,43	0
2	MLY	B	624	11/12	0.96	0.14	-	27,31,57,58	0
2	MLY	B	946	11/12	0.95	0.11	-	26,34,46,48	0
2	MLY	B	642	11/12	0.96	0.10	-	27,28,39,43	0
2	MLY	B	986	11/12	0.92	0.17	-	35,43,70,70	0
1	MLY	A	750	11/12	0.94	0.12	-	27,41,61,66	0
2	MLZ	B	769	10/11	0.93	0.11	-	30,41,50,52	0
1	MLY	A	911	11/12	0.96	0.14	-	20,22,42,44	0
1	MLY	A	937	11/12	0.94	0.13	-	24,31,62,65	0
1	MLY	A	154	11/12	0.98	0.08	-	23,28,61,63	0
2	MLY	B	488	11/12	0.98	0.09	-	27,30,42,43	0
2	MLY	B	956	11/12	0.96	0.12	-	25,34,64,67	0
1	MLY	A	794	11/12	0.94	0.14	-	38,52,69,69	0
1	MLY	A	363	11/12	0.95	0.11	-	24,31,59,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MLZ	B	937	10/11	0.96	0.12	-	23,28,64,67	0
2	MLY	B	287	11/12	0.95	0.12	-	31,35,50,52	0
1	CAS	A	477	9/10	0.89	0.11	-	40,44,77,147	0
2	MLY	B	499	11/12	0.94	0.14	-	28,33,64,67	0
2	MLY	B	943	11/12	0.96	0.10	-	24,31,56,58	0
1	MLY	A	854	11/12	0.96	0.09	-	20,29,46,46	0
2	MLY	B	199	11/12	0.95	0.14	-	22,31,62,62	0
1	MLZ	A	550	10/11	0.96	0.11	-	28,35,45,46	0
1	MLY	A	466	11/12	0.97	0.10	-	28,32,50,51	0
1	MLY	A	278	11/12	0.97	0.08	-	22,35,41,41	0
1	MLY	A	66	11/12	0.96	0.11	-	25,36,54,55	0
2	CAS	B	171	9/10	0.90	0.17	-	24,32,90,102	0
1	MLY	A	642	11/12	0.98	0.09	-	22,24,42,43	0
1	MLY	A	287	11/12	0.98	0.10	-	22,26,34,34	0
1	MLY	A	946	11/12	0.97	0.12	-	22,33,57,60	0
2	MLZ	B	884	10/11	0.97	0.08	-	22,27,53,54	0
1	MLY	A	704	11/12	0.93	0.12	-	25,29,50,52	0
2	MLY	B	911	11/12	0.97	0.10	-	20,28,54,55	0
2	MLY	B	431	11/12	0.98	0.14	-	25,32,60,62	0
1	MLZ	A	437	10/11	0.98	0.08	-	28,32,56,56	0
2	MLY	B	704	11/12	0.97	0.10	-	24,31,59,60	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
5	ACT	B	1114	4/4	0.85	0.31	15.25	61,70,71,74	0
4	GOL	B	1112	6/6	0.59	0.26	13.04	84,89,91,92	0
5	ACT	A	1104	4/4	0.93	0.29	10.06	49,57,57,58	0
5	ACT	B	1109	4/4	0.78	0.20	6.48	91,91,92,93	0
5	ACT	B	1110	4/4	0.97	0.16	5.70	8,31,41,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	ACT	A	1105	4/4	0.54	0.33	5.68	58,64,65,72	0
5	ACT	B	1113	4/4	0.68	0.18	5.62	49,56,60,62	0
5	ACT	A	1108	4/4	0.85	0.23	5.08	34,35,35,36	0
5	ACT	B	1107	4/4	0.89	0.19	4.52	69,71,73,77	0
4	GOL	B	1102	6/6	0.97	0.17	3.96	11,48,54,64	0
5	ACT	B	1105	4/4	0.71	0.17	3.69	49,59,60,66	0
5	ACT	B	1106	4/4	0.93	0.13	3.07	69,71,71,74	0
5	ACT	A	1106	4/4	0.27	0.22	2.38	58,65,67,67	0
4	GOL	A	1107	6/6	0.87	0.20	2.35	40,53,63,68	0
5	ACT	B	1111	4/4	0.73	0.15	1.85	72,76,77,81	0
5	ACT	B	1108	4/4	0.97	0.15	0.28	38,40,43,47	0
3	ZN	A	1101	1/1	0.95	0.11	0.10	28,28,28,28	0
4	GOL	B	1104	6/6	0.94	0.09	0.02	35,47,50,52	0
4	GOL	A	1103	6/6	0.93	0.08	-0.41	40,41,50,56	0
4	GOL	B	1103	6/6	0.98	0.08	-0.94	23,32,36,39	0
4	GOL	A	1102	6/6	0.97	0.08	-1.46	23,33,34,35	0
3	ZN	B	1101	1/1	0.99	0.09	-2.27	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.