



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:06 AM GMT

PDB ID : 3PII
Title : Crystal structure of Mutant of ht- Alcohol Dehydrogenase with substrate analogue butyramide
Authors : Dong, M.; Bahnson, B.J.
Deposited on : 2010-11-06
Resolution : 2.90 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

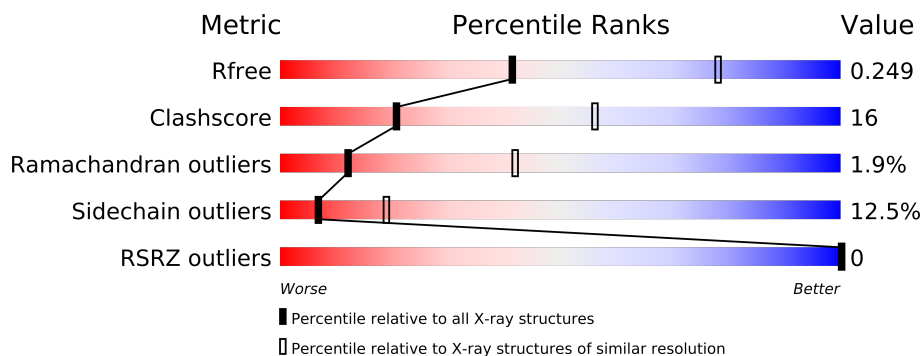
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.90 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}	66092	1053 (2.90-2.90)
Clashscore	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (2.90-2.90)
RSRZ outliers	66119	1054 (2.90-2.90)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	339	
1	B	339	
1	C	339	
1	D	339	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	BMD	A	600	-	X
3	BMD	B	500	-	X
3	BMD	C	700	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10211 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Alcohol dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2539	1630	427	468	14			
1	C	337	Total	C	N	O	S	0	0	0
			2539	1630	427	468	14			
1	B	337	Total	C	N	O	S	0	0	0
			2539	1630	427	468	14			
1	D	337	Total	C	N	O	S	0	0	0
			2539	1630	427	468	14			

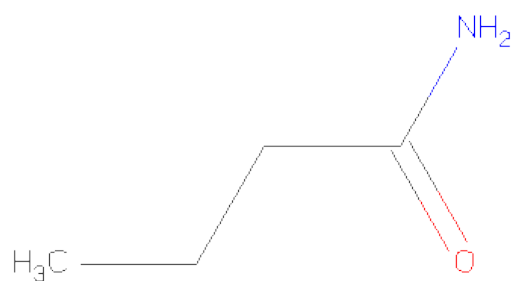
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	176	PHE	LEU	ENGINEERED MUTATION	UNP P42328
C	176	PHE	LEU	ENGINEERED MUTATION	UNP P42328
B	176	PHE	LEU	ENGINEERED MUTATION	UNP P42328
D	176	PHE	LEU	ENGINEERED MUTATION	UNP P42328

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

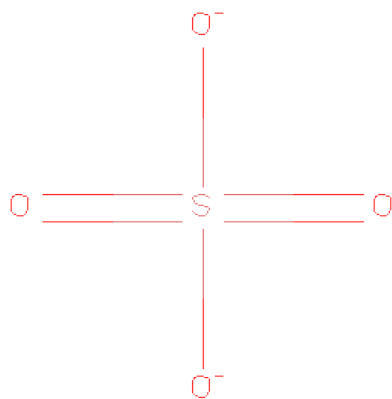
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		

- Molecule 3 is BUTYRAMIDE (three-letter code: BMD) (formula: C₄H₉NO).



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

- Molecule 4 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

- Molecule 5 is water.

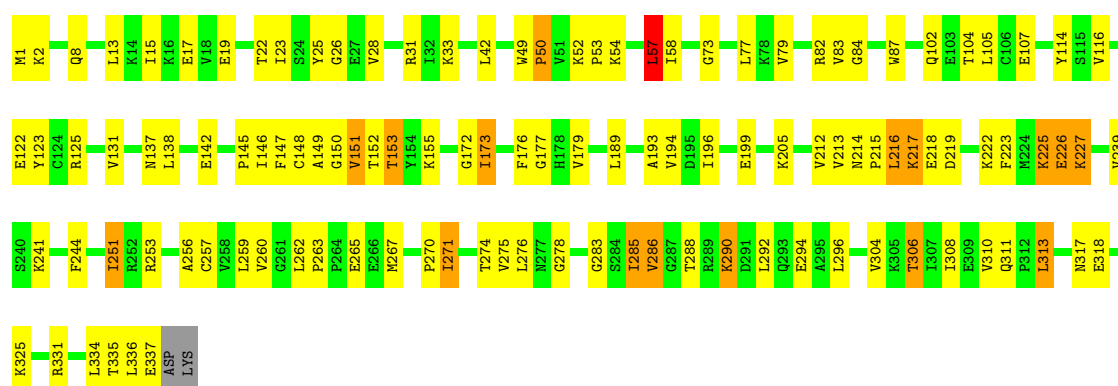
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O 1 1	0	0
5	C	1	Total O 1 1	0	0
5	B	1	Total O 1 1	0	0

3 Residue-property plots

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.

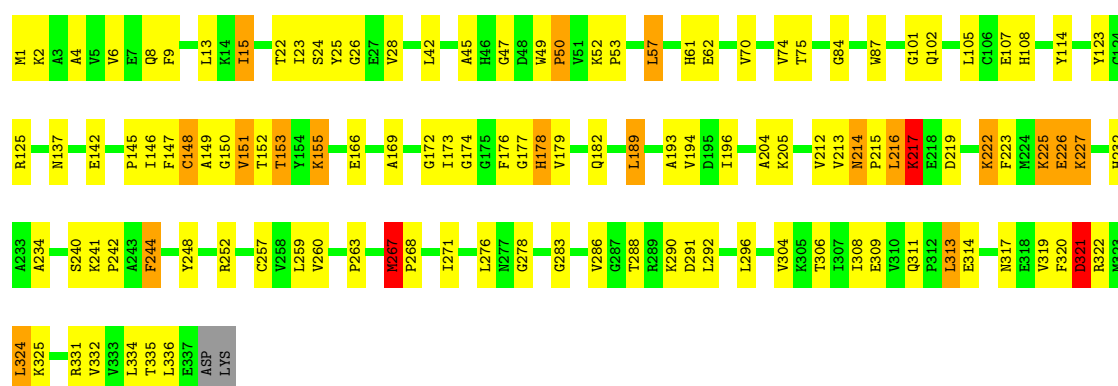
- Molecule 1: Alcohol dehydrogenase

Chain A: 



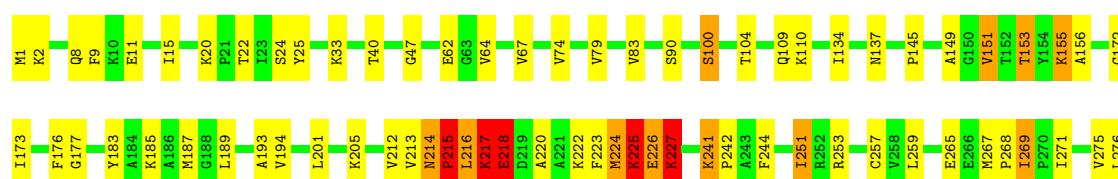
- Molecule 1: Alcohol dehydrogenase

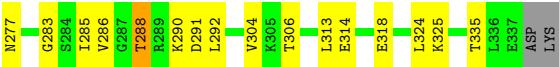
Chain C: 



- Molecule 1: Alcohol dehydrogenase

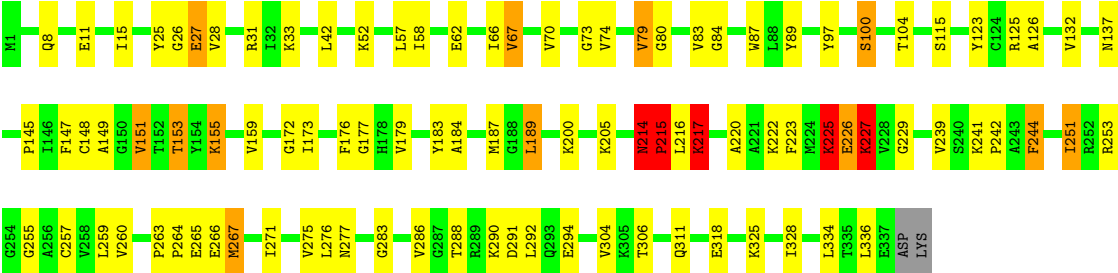
Chain B: 





● Molecule 1: Alcohol dehydrogenase

Chain D: A horizontal bar chart for Chain D. The bar is mostly green, with a yellow segment towards the right end and a small orange segment at the far right.



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	145.65Å 145.65Å 135.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.61 – 2.90 49.61 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.61-2.90) 99.8 (49.61-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.41 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.206 , 0.286 0.198 , 0.249	Depositor DCC
R_{free} test set	1799 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	46.1	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 5.3	EDS
Estimated twinning fraction	0.488 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	0 of 36059 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10211	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, SO4, BMD

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	0.75	0/2591	0.81	3/3513 (0.1%)
1	B	0.72	0/2591	0.83	1/3513 (0.0%)
1	C	0.78	2/2591 (0.1%)	0.85	6/3513 (0.2%)
1	D	0.71	0/2591	0.79	3/3513 (0.1%)
All	All	0.74	2/10364 (0.0%)	0.82	13/14052 (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
1	B	0	8
1	C	0	5
1	D	0	4
All	All	0	18

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	257	CYS	CB-SG	-5.74	1.72	1.81
1	C	309	GLU	CG-CD	5.09	1.59	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	226	GLU	C-N-CA	5.98	136.64	121.70
1	D	214	ASN	N-CA-C	5.87	126.85	111.00
1	D	215	PRO	N-CA-C	-5.81	97.01	112.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	321	ASP	N-CA-C	5.73	126.47	111.00
1	B	225	LYS	CB-CA-C	5.61	121.62	110.40
1	C	252	ARG	NE-CZ-NH1	5.47	123.03	120.30
1	C	215	PRO	N-CA-C	-5.21	98.55	112.10
1	C	189	LEU	CA-CB-CG	5.16	127.17	115.30
1	C	57	LEU	CA-CB-CG	5.14	127.13	115.30
1	A	57	LEU	CA-CB-CG	5.12	127.07	115.30
1	A	173	ILE	CG1-CB-CG2	-5.10	100.19	111.40
1	A	226	GLU	C-N-CA	5.07	134.36	121.70
1	C	226	GLU	C-N-CA	5.01	134.23	121.70

There are no chirality outliers.

All (18) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	225	LYS	Peptide
1	B	214	ASN	Peptide
1	B	215	PRO	Peptide
1	B	217	LYS	Peptide
1	B	218	GLU	Peptide
1	B	224	MET	Peptide
1	B	225	LYS	Peptide
1	B	226	GLU	Mainchain,Peptide
1	C	148	CYS	Peptide
1	C	214	ASN	Peptide
1	C	217	LYS	Peptide
1	C	225	LYS	Peptide
1	C	321	ASP	Peptide
1	D	214	ASN	Peptide
1	D	215	PRO	Peptide
1	D	217	LYS	Peptide
1	D	225	LYS	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2539	0	2586	82	0
1	B	2539	0	2586	95	0
1	C	2539	0	2586	93	0
1	D	2539	0	2586	80	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	6	0	9	2	0
3	B	6	0	9	3	0
3	C	6	0	9	0	0
3	D	6	0	9	2	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
4	D	5	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
All	All	10211	0	10380	322	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 16.

All (322) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:215:PRO:HD2	1:B:216:LEU:CD2	1.56	1.34
1:B:215:PRO:CD	1:B:216:LEU:HD22	1.54	1.33
1:B:214:ASN:OD1	1:B:216:LEU:HD21	1.16	1.29
1:B:226:GLU:H	1:B:227:LYS:CB	1.48	1.26
1:B:215:PRO:CD	1:B:216:LEU:CD2	2.11	1.20
1:C:216:LEU:CB	1:C:217:LYS:HB2	1.69	1.19
1:B:216:LEU:O	1:B:217:LYS:HG2	1.39	1.18
1:D:239:VAL:HG12	1:D:267:MET:CE	1.77	1.14
1:B:214:ASN:OD1	1:B:216:LEU:CD2	1.98	1.10
1:C:216:LEU:HB3	1:C:217:LYS:HB2	1.19	1.08
1:D:288:THR:HG22	1:D:290:LYS:H	1.14	1.08
1:B:216:LEU:O	1:B:216:LEU:HD23	1.55	1.04
1:B:226:GLU:H	1:B:227:LYS:HB2	1.19	1.03
1:B:226:GLU:N	1:B:227:LYS:CB	2.22	1.02
1:B:223:PHE:O	1:B:227:LYS:HG2	1.61	0.99
1:B:226:GLU:H	1:B:227:LYS:HB3	1.22	0.99
1:B:217:LYS:O	1:B:217:LYS:HG3	1.63	0.98

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:196:ILE:O	1:A:214:ASN:ND2	1.98	0.96
1:D:214:ASN:O	1:D:216:LEU:HB3	1.66	0.94
1:C:216:LEU:CB	1:C:217:LYS:CB	2.46	0.94
1:B:215:PRO:N	1:B:216:LEU:HD23	1.84	0.93
1:B:216:LEU:O	1:B:217:LYS:CG	2.15	0.93
1:B:226:GLU:N	1:B:227:LYS:HB2	1.84	0.92
1:D:288:THR:HG22	1:D:290:LYS:N	1.88	0.89
1:D:239:VAL:HG12	1:D:267:MET:HE2	1.55	0.89
1:D:216:LEU:O	1:D:217:LYS:HB3	1.73	0.88
1:A:226:GLU:HB3	1:A:227:LYS:HB2	1.55	0.88
1:B:215:PRO:N	1:B:216:LEU:CD2	2.37	0.86
1:C:150:GLY:O	1:C:153:THR:HG22	1.76	0.84
1:C:288:THR:HG22	1:C:290:LYS:H	1.42	0.83
1:D:214:ASN:OD1	1:D:214:ASN:O	1.97	0.82
1:C:13:LEU:H	1:C:317:ASN:HD21	1.27	0.79
1:B:216:LEU:O	1:B:216:LEU:CD2	2.30	0.79
1:A:288:THR:HG22	1:A:290:LYS:H	1.48	0.79
1:C:263:PRO:HD2	1:B:271:ILE:HG21	1.65	0.79
1:B:226:GLU:N	1:B:227:LYS:HB3	1.92	0.78
1:B:223:PHE:O	1:B:227:LYS:CG	2.30	0.78
1:C:216:LEU:HB2	1:C:217:LYS:CB	2.12	0.78
1:B:40:THR:HG21	3:B:500:BMD:H11	1.64	0.77
1:C:263:PRO:HD2	1:B:271:ILE:CG2	2.15	0.77
1:B:288:THR:HG22	1:B:290:LYS:N	1.98	0.77
1:A:222:LYS:O	1:A:226:GLU:HB2	1.87	0.74
1:B:201:LEU:HD11	1:B:214:ASN:HB2	1.69	0.73
1:D:149:ALA:O	1:D:153:THR:HB	1.89	0.72
1:B:151:VAL:HG13	1:B:292:LEU:HA	1.72	0.72
1:C:102:GLN:HB3	1:C:105:LEU:HD12	1.72	0.72
1:C:13:LEU:N	1:C:317:ASN:HD21	1.87	0.71
1:B:149:ALA:O	1:B:153:THR:HB	1.91	0.71
1:B:288:THR:HG22	1:B:290:LYS:H	1.55	0.70
1:A:146:ILE:HG21	1:A:296:LEU:HD23	1.72	0.70
1:D:239:VAL:HG12	1:D:267:MET:HE1	1.69	0.70
1:A:13:LEU:H	1:A:317:ASN:HD21	1.39	0.70
1:C:271:ILE:HD11	1:B:267:MET:HB2	1.74	0.69
1:C:267:MET:HB3	1:C:268:PRO:CA	2.22	0.69
1:A:263:PRO:HD2	1:D:271:ILE:CG2	2.23	0.69
1:B:226:GLU:CA	1:B:227:LYS:HB2	2.22	0.68
1:A:193:ALA:HB3	1:A:212:VAL:HG12	1.74	0.68
1:B:214:ASN:O	1:B:217:LYS:HG2	1.94	0.68
1:D:67:VAL:HG12	1:D:79:VAL:HA	1.77	0.67

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:13:LEU:N	1:A:317:ASN:HD21	1.93	0.67
1:C:151:VAL:HG13	1:C:292:LEU:HD13	1.77	0.67
1:C:244:PHE:HD1	1:C:244:PHE:C	1.98	0.66
1:C:42:LEU:HD11	1:C:324:LEU:HD12	1.77	0.66
1:B:216:LEU:O	1:B:217:LYS:CB	2.43	0.66
1:B:2:LYS:NZ	1:B:314:GLU:OE2	2.29	0.66
1:C:216:LEU:HB2	1:C:217:LYS:HB2	1.65	0.66
1:C:244:PHE:CD1	1:C:244:PHE:C	2.70	0.65
1:A:219:ASP:OD2	1:A:222:LYS:HB2	1.96	0.65
1:D:251:ILE:HD13	1:D:257:CYS:HB2	1.79	0.65
1:D:151:VAL:HG22	1:D:292:LEU:HD13	1.77	0.65
1:A:150:GLY:O	1:A:153:THR:HG22	1.96	0.64
1:C:173:ILE:O	1:C:173:ILE:HG22	1.97	0.64
1:C:149:ALA:O	1:C:153:THR:HB	1.98	0.64
1:A:267:MET:HB3	1:D:271:ILE:HD13	1.80	0.64
1:B:216:LEU:C	1:B:216:LEU:HD23	2.15	0.64
1:A:153:THR:HG21	1:A:179:VAL:HG23	1.79	0.64
1:A:102:GLN:HB3	1:A:105:LEU:HD12	1.79	0.63
1:B:215:PRO:HD2	1:B:216:LEU:HD22	0.70	0.63
1:C:146:ILE:HG21	1:C:296:LEU:HD23	1.80	0.63
1:A:263:PRO:HD2	1:D:271:ILE:HG21	1.80	0.63
1:B:251:ILE:HD13	1:B:257:CYS:HB2	1.81	0.62
1:B:251:ILE:CD1	1:B:257:CYS:HB2	2.29	0.62
1:C:222:LYS:O	1:C:226:GLU:HB2	1.99	0.62
1:A:142:GLU:O	1:A:145:PRO:HD2	2.00	0.62
1:A:271:ILE:CD1	1:D:267:MET:HB2	2.30	0.62
1:C:267:MET:HG2	1:B:271:ILE:HD11	1.81	0.62
1:D:239:VAL:CG1	1:D:267:MET:HE2	2.30	0.62
1:D:62:GLU:OE1	1:D:145:PRO:HA	2.00	0.62
1:C:155:LYS:HG2	1:C:291:ASP:OD2	2.00	0.61
1:B:193:ALA:HB3	1:B:212:VAL:HG12	1.82	0.61
1:A:218:GLU:HG2	1:A:219:ASP:N	2.14	0.61
1:B:214:ASN:C	1:B:216:LEU:HD23	2.20	0.61
1:B:8:GLN:HB2	1:B:11:GLU:HG3	1.82	0.61
1:C:216:LEU:HB2	1:C:217:LYS:HB3	1.83	0.60
1:A:278:GLY:HA2	1:D:283:GLY:O	2.02	0.60
1:B:288:THR:CG2	1:B:290:LYS:H	2.14	0.60
1:A:146:ILE:CG2	1:A:296:LEU:HD23	2.31	0.60
1:B:244:PHE:HZ	1:B:269:ILE:HG22	1.66	0.60
1:A:263:PRO:HG2	1:D:271:ILE:HG21	1.83	0.59
1:A:23:ILE:HG22	1:A:125:ARG:HG3	1.84	0.59
1:C:13:LEU:H	1:C:317:ASN:ND2	2.00	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:216:LEU:O	1:D:217:LYS:CB	2.49	0.59
1:A:271:ILE:HD11	1:D:267:MET:HB2	1.84	0.59
1:A:2:LYS:HG2	1:A:122:GLU:OE2	2.03	0.58
1:C:142:GLU:O	1:C:145:PRO:HD2	2.04	0.58
1:C:267:MET:HB3	1:C:268:PRO:HA	1.85	0.58
1:C:4:ALA:HB2	1:C:313:LEU:HD21	1.86	0.58
1:C:268:PRO:HA	1:B:268:PRO:HA	1.85	0.58
1:C:194:VAL:HG22	1:C:213:VAL:HG12	1.84	0.57
1:D:153:THR:HG21	1:D:179:VAL:HG23	1.86	0.57
1:D:173:ILE:O	1:D:200:LYS:HE2	2.04	0.57
1:C:319:VAL:HG11	1:C:332:VAL:HG11	1.87	0.56
1:C:172:GLY:O	1:C:177:GLY:HA3	2.04	0.56
1:C:153:THR:HG21	1:C:179:VAL:HG23	1.88	0.56
1:C:62:GLU:OE1	1:C:145:PRO:HA	2.05	0.56
1:D:155:LYS:HD2	1:D:159:VAL:HG23	1.88	0.56
1:D:15:ILE:H	1:D:15:ILE:HD12	1.69	0.56
1:B:241:LYS:HG2	1:B:242:PRO:HD3	1.87	0.56
1:D:28:VAL:HG12	1:D:70:VAL:HA	1.87	0.56
1:D:57:LEU:HD22	1:D:58:ILE:N	2.21	0.55
1:B:215:PRO:CD	1:B:216:LEU:HD21	2.29	0.55
1:C:214:ASN:C	1:C:216:LEU:H	2.01	0.55
1:D:244:PHE:HD1	1:D:244:PHE:C	2.09	0.55
1:C:226:GLU:H	1:C:227:LYS:HB3	1.71	0.55
1:C:23:ILE:O	1:C:23:ILE:HG13	2.07	0.55
1:A:57:LEU:HD22	1:A:58:ILE:N	2.22	0.55
1:A:263:PRO:CG	1:D:271:ILE:HG21	2.37	0.54
1:A:251:ILE:HD11	1:A:256:ALA:O	2.07	0.54
1:C:244:PHE:O	1:C:244:PHE:HD1	1.90	0.54
1:C:182:GLN:CB	1:C:304:VAL:HG12	2.37	0.54
1:B:217:LYS:O	1:B:217:LYS:CG	2.42	0.54
1:B:216:LEU:O	1:B:216:LEU:CG	2.55	0.54
1:A:311:GLN:O	1:A:334:LEU:HA	2.08	0.54
1:D:244:PHE:CD1	1:D:244:PHE:C	2.81	0.54
1:D:244:PHE:O	1:D:244:PHE:HD1	1.90	0.53
1:B:15:ILE:HD12	1:B:15:ILE:H	1.71	0.53
1:C:263:PRO:CD	1:B:271:ILE:HG21	2.35	0.53
1:A:149:ALA:HA	1:A:176:PHE:CE2	2.44	0.53
1:C:226:GLU:HB3	1:C:227:LYS:HB2	1.91	0.53
1:C:149:ALA:HB1	1:C:176:PHE:CD2	2.43	0.53
1:C:288:THR:HG22	1:C:290:LYS:N	2.19	0.52
1:D:151:VAL:CG2	1:D:292:LEU:HD13	2.39	0.52
1:D:184:ALA:O	1:D:189:LEU:HB2	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:223:PHE:CE2	1:A:227:LYS:HE3	2.45	0.52
1:C:214:ASN:O	1:C:216:LEU:N	2.32	0.52
1:C:174:GLY:O	1:C:178:HIS:HB3	2.10	0.52
1:A:263:PRO:CD	1:D:271:ILE:HG21	2.40	0.52
1:A:173:ILE:CD1	1:A:193:ALA:HB1	2.40	0.52
1:D:251:ILE:HG23	1:D:255:GLY:HA3	1.92	0.52
1:D:226:GLU:H	1:D:227:LYS:HB3	1.75	0.52
1:B:223:PHE:O	1:B:227:LYS:CB	2.58	0.52
1:C:214:ASN:O	1:C:217:LYS:HB3	2.09	0.52
1:A:179:VAL:HG11	1:A:306:THR:HG23	1.92	0.51
1:C:149:ALA:HB1	1:C:176:PHE:HD2	1.74	0.51
1:D:67:VAL:HG12	1:D:67:VAL:O	2.10	0.51
1:C:263:PRO:HD2	1:B:271:ILE:HG22	1.93	0.51
1:B:183:TYR:O	1:B:187:MET:HG3	2.10	0.51
1:B:214:ASN:O	1:B:217:LYS:CG	2.59	0.51
1:A:1:MET:SD	1:A:123:TYR:HB2	2.50	0.51
1:D:26:GLY:HA2	1:D:74:VAL:HG12	1.92	0.51
1:A:25:TYR:CE2	1:C:25:TYR:CE2	2.99	0.51
1:C:283:GLY:HA3	1:B:275:VAL:HA	1.92	0.51
1:D:151:VAL:HG13	1:D:292:LEU:HA	1.93	0.51
1:B:220:ALA:O	1:B:224:MET:HB2	2.11	0.50
1:C:219:ASP:CG	1:C:222:LYS:HB2	2.31	0.50
1:A:283:GLY:HA3	1:D:275:VAL:HA	1.92	0.50
1:B:214:ASN:OD1	1:B:214:ASN:C	2.49	0.50
1:A:33:LYS:NZ	1:A:337:GLU:HG3	2.27	0.50
1:A:173:ILE:HD11	1:A:193:ALA:HB1	1.93	0.50
1:C:146:ILE:CG2	1:C:296:LEU:HD23	2.42	0.50
1:B:217:LYS:C	1:B:218:GLU:O	2.50	0.50
1:C:216:LEU:CA	1:C:217:LYS:CB	2.90	0.50
1:D:172:GLY:O	1:D:177:GLY:HA3	2.11	0.49
1:C:244:PHE:CE1	1:C:248:TYR:HB2	2.47	0.49
1:C:28:VAL:HG12	1:C:70:VAL:HG12	1.94	0.49
1:D:264:PRO:HG2	1:D:265:GLU:H	1.77	0.49
1:A:285:ILE:HG23	1:A:286:VAL:HG22	1.94	0.49
1:C:214:ASN:C	1:C:214:ASN:OD1	2.50	0.49
1:B:244:PHE:CZ	1:B:269:ILE:HG22	2.47	0.49
1:C:61:HIS:HB2	1:C:87:TRP:HA	1.94	0.49
1:C:23:ILE:HG22	1:C:125:ARG:HG3	1.94	0.49
1:C:182:GLN:HB2	1:C:304:VAL:HG12	1.95	0.49
1:C:25:TYR:CD2	1:C:26:GLY:N	2.81	0.49
1:B:62:GLU:OE1	1:B:145:PRO:HA	2.12	0.49
1:B:217:LYS:O	1:B:218:GLU:HG2	2.13	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:251:ILE:CD1	1:D:257:CYS:HB2	2.41	0.49
1:A:290:LYS:HD3	1:A:294:GLU:OE2	2.13	0.48
1:A:244:PHE:C	1:A:244:PHE:CD1	2.87	0.48
1:C:84:GLY:HA3	1:C:147:PHE:CE2	2.48	0.48
1:D:288:THR:CG2	1:D:290:LYS:H	2.04	0.48
1:B:67:VAL:HG12	1:B:79:VAL:HA	1.96	0.48
1:D:241:LYS:HB2	1:D:242:PRO:HD3	1.94	0.48
1:A:33:LYS:HZ2	1:A:337:GLU:HG3	1.79	0.48
1:D:27:GLU:OE1	1:D:125:ARG:HD2	2.13	0.48
1:C:15:ILE:HG21	1:C:314:GLU:HG3	1.94	0.48
1:D:8:GLN:HB2	1:D:11:GLU:HG3	1.95	0.48
1:A:172:GLY:O	1:A:177:GLY:HA3	2.13	0.48
1:C:271:ILE:CD1	1:B:267:MET:HB2	2.42	0.47
1:A:84:GLY:HA3	1:A:147:PHE:CE2	2.48	0.47
1:A:77:LEU:HD11	1:A:131:VAL:HG22	1.96	0.47
1:C:278:GLY:HA2	1:B:283:GLY:O	2.14	0.47
1:C:4:ALA:HB2	1:C:313:LEU:CD2	2.44	0.47
1:A:313:LEU:HD22	1:A:313:LEU:O	2.15	0.47
1:D:215:PRO:HB3	1:D:220:ALA:HB2	1.97	0.47
1:D:227:LYS:HA	1:D:227:LYS:HD3	1.62	0.47
1:D:87:TRP:CE2	3:D:400:BMD:H21	2.49	0.47
1:B:225:LYS:HA	1:B:225:LYS:HD3	1.77	0.47
1:D:225:LYS:O	1:D:229:GLY:HA2	2.15	0.46
1:B:214:ASN:O	1:B:216:LEU:O	2.33	0.46
1:A:239:VAL:HB	1:A:263:PRO:HG3	1.97	0.46
1:B:40:THR:CG2	3:B:500:BMD:H11	2.38	0.46
1:C:166:GLU:HG2	1:C:232:HIS:CE1	2.50	0.46
1:A:49:TRP:O	1:A:50:PRO:C	2.52	0.46
1:C:241:LYS:HG2	1:C:241:LYS:H	1.56	0.46
1:D:66:ILE:HG23	1:D:80:GLY:HA2	1.96	0.46
1:A:104:THR:HG23	1:A:105:LEU:HG	1.97	0.46
1:C:1:MET:SD	1:C:123:TYR:HB2	2.56	0.46
1:A:23:ILE:HG13	1:A:23:ILE:O	2.16	0.46
1:D:89:TYR:CE1	1:D:126:ALA:HA	2.51	0.46
1:B:151:VAL:CG1	1:B:292:LEU:HA	2.44	0.46
1:C:173:ILE:HG12	1:C:204:ALA:HB2	1.98	0.46
1:D:226:GLU:H	1:D:227:LYS:CB	2.29	0.45
1:D:225:LYS:HA	1:D:225:LYS:HD2	1.77	0.45
1:B:194:VAL:HG22	1:B:213:VAL:HG12	1.99	0.45
1:A:173:ILE:HG22	1:A:173:ILE:O	2.15	0.45
1:A:214:ASN:OD1	1:A:216:LEU:N	2.49	0.45
1:D:215:PRO:HG2	1:D:216:LEU:HB2	1.99	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:222:LYS:O	1:B:226:GLU:HB2	2.16	0.45
1:A:288:THR:HG22	1:A:290:LYS:N	2.23	0.45
1:B:213:VAL:HG22	1:B:214:ASN:H	1.82	0.45
1:B:15:ILE:N	1:B:15:ILE:HD12	2.31	0.45
1:A:194:VAL:HG22	1:A:213:VAL:HG13	1.99	0.45
1:D:288:THR:HB	1:D:291:ASP:H	1.82	0.45
1:B:226:GLU:CB	1:B:227:LYS:HB2	2.47	0.45
1:C:2:LYS:HE3	1:C:314:GLU:OE2	2.17	0.45
1:A:84:GLY:HA3	1:A:147:PHE:CZ	2.52	0.44
1:D:311:GLN:O	1:D:334:LEU:HA	2.17	0.44
1:A:270:PRO:O	1:A:274:THR:OG1	2.23	0.44
1:B:227:LYS:HD3	1:B:227:LYS:HA	1.42	0.44
1:A:145:PRO:HB2	1:A:306:THR:HG21	1.99	0.44
1:A:25:TYR:CD2	1:A:26:GLY:N	2.86	0.44
1:B:214:ASN:C	1:B:216:LEU:CD2	2.84	0.44
1:B:110:LYS:HD3	1:B:110:LYS:HA	1.86	0.44
1:A:270:PRO:HA	1:D:266:GLU:HG2	1.98	0.44
1:B:201:LEU:CD1	1:B:214:ASN:HB2	2.45	0.44
1:C:244:PHE:O	1:C:244:PHE:CD1	2.70	0.44
1:C:173:ILE:CG2	1:C:173:ILE:O	2.66	0.44
1:C:193:ALA:HB3	1:C:212:VAL:HG12	2.00	0.44
1:A:267:MET:HB3	1:D:271:ILE:CD1	2.47	0.44
1:C:151:VAL:O	1:C:152:THR:C	2.55	0.44
1:A:148:CYS:SG	3:A:600:BMD:NA	2.82	0.44
1:D:84:GLY:HA3	1:D:147:PHE:CE2	2.53	0.43
1:B:90:SER:O	1:B:109:GLN:HA	2.18	0.43
1:A:214:ASN:HA	1:A:215:PRO:HD3	1.86	0.43
1:D:222:LYS:O	1:D:226:GLU:HB2	2.18	0.43
1:D:253:ARG:HG2	1:D:277:ASN:HB3	2.00	0.43
1:A:151:VAL:O	1:A:152:THR:C	2.57	0.43
1:C:6:VAL:HG11	1:C:45:ALA:HA	2.01	0.43
1:B:216:LEU:N	1:B:216:LEU:CD2	2.81	0.43
1:A:271:ILE:HD13	1:D:267:MET:HB2	1.99	0.43
1:A:310:VAL:O	1:A:311:GLN:HG2	2.19	0.43
1:D:223:PHE:O	1:D:227:LYS:CB	2.65	0.43
1:A:25:TYR:CZ	1:A:73:GLY:HA3	2.53	0.43
1:B:241:LYS:HB3	1:B:267:MET:HG2	2.00	0.43
1:B:288:THR:HG22	1:B:291:ASP:H	1.84	0.43
1:B:149:ALA:HA	1:B:176:PHE:CE2	2.54	0.43
1:C:26:GLY:HA2	1:C:74:VAL:HG12	2.01	0.43
1:B:67:VAL:HG12	1:B:67:VAL:O	2.18	0.43
1:A:53:PRO:HD3	1:A:114:TYR:CZ	2.54	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:169:ALA:O	1:C:234:ALA:HA	2.18	0.43
1:D:31:ARG:HB2	1:D:123:TYR:CE1	2.53	0.43
1:B:172:GLY:O	1:B:177:GLY:HA3	2.19	0.43
1:D:132:VAL:HG21	1:D:292:LEU:HD23	2.01	0.43
1:C:196:ILE:O	1:C:214:ASN:ND2	2.49	0.43
1:D:216:LEU:HG	1:D:216:LEU:O	2.18	0.43
1:C:70:VAL:HB	1:C:74:VAL:CG2	2.49	0.43
1:A:241:LYS:HE2	1:A:265:GLU:HG3	2.01	0.43
1:A:149:ALA:O	1:A:153:THR:HB	2.18	0.42
1:C:223:PHE:O	1:C:227:LYS:HB3	2.19	0.42
1:A:214:ASN:C	1:A:216:LEU:H	2.22	0.42
1:C:311:GLN:O	1:C:334:LEU:HA	2.19	0.42
1:A:82:ARG:NH2	1:A:138:LEU:O	2.42	0.42
1:C:226:GLU:H	1:C:227:LYS:CB	2.32	0.42
1:A:275:VAL:HA	1:D:283:GLY:HA3	2.01	0.42
1:B:253:ARG:HG2	1:B:277:ASN:HB3	2.00	0.42
1:C:240:SER:OG	1:C:242:PRO:HD2	2.20	0.42
1:B:213:VAL:HG22	1:B:214:ASN:N	2.34	0.42
1:A:216:LEU:CB	1:A:217:LYS:HB3	2.50	0.42
1:A:87:TRP:CZ2	3:A:600:BMD:H32	2.55	0.42
1:A:262:LEU:N	1:A:263:PRO:HD3	2.34	0.42
1:C:223:PHE:O	1:C:227:LYS:CB	2.67	0.42
1:A:227:LYS:HD3	1:A:227:LYS:HA	1.88	0.42
1:B:40:THR:HG21	3:B:500:BMD:C1	2.42	0.41
1:A:271:ILE:CG2	1:D:263:PRO:HD2	2.49	0.41
1:A:267:MET:CB	1:D:271:ILE:HD13	2.47	0.41
1:A:223:PHE:O	1:A:226:GLU:HB3	2.21	0.41
1:B:9:PHE:CD1	1:B:47:GLY:HA3	2.56	0.41
1:D:25:TYR:CE2	1:D:73:GLY:HA3	2.56	0.41
1:D:290:LYS:O	1:D:294:GLU:HG3	2.21	0.41
1:A:214:ASN:OD1	1:A:216:LEU:HB2	2.21	0.41
1:C:151:VAL:CG1	1:C:292:LEU:HD13	2.48	0.41
1:B:267:MET:HA	1:B:268:PRO:HD3	1.77	0.41
1:D:147:PHE:CD2	1:D:292:LEU:HD21	2.56	0.41
1:C:223:PHE:CE2	1:C:227:LYS:HE3	2.56	0.41
1:C:84:GLY:HA3	1:C:147:PHE:CZ	2.56	0.41
1:D:148:CYS:SG	3:D:400:BMD:OA	2.79	0.41
1:A:253:ARG:O	1:D:97:TYR:HE1	2.04	0.41
1:B:24:SER:HB3	1:B:25:TYR:H	1.52	0.41
1:B:185:LYS:HA	1:B:185:LYS:HD2	1.90	0.41
1:B:155:LYS:O	1:B:156:ALA:C	2.59	0.41
1:B:1:MET:HB3	1:B:20:LYS:HB2	2.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:183:TYR:O	1:D:187:MET:HG3	2.21	0.41
1:C:53:PRO:HD3	1:C:114:TYR:CZ	2.56	0.40
1:D:216:LEU:O	1:D:216:LEU:CG	2.68	0.40
1:B:64:VAL:HB	1:B:134:ILE:HD11	2.03	0.40
1:C:49:TRP:O	1:C:50:PRO:C	2.59	0.40
1:D:149:ALA:HA	1:D:176:PHE:CE2	2.56	0.40
1:C:182:GLN:HB3	1:C:304:VAL:HG12	2.04	0.40
1:B:173:ILE:HA	1:B:177:GLY:HA3	2.02	0.40
1:C:9:PHE:CD1	1:C:47:GLY:HA3	2.56	0.40
1:A:151:VAL:HG13	1:A:292:LEU:HD13	2.04	0.40
1:C:321:ASP:H	1:C:322:ARG:HB2	1.87	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	335/339 (99%)	299 (89%)	31 (9%)	5 (2%)	15	50
1	B	335/339 (99%)	301 (90%)	27 (8%)	7 (2%)	11	39
1	C	335/339 (99%)	296 (88%)	32 (10%)	7 (2%)	11	39
1	D	335/339 (99%)	303 (90%)	26 (8%)	6 (2%)	13	44
All	All	1340/1356 (99%)	1199 (90%)	116 (9%)	25 (2%)	12	42

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	227	LYS
1	C	227	LYS
1	C	267	MET
1	B	215	PRO
1	B	217	LYS
1	B	218	GLU
1	B	225	LYS

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Mol	Chain	Res	Type
1	B	227	LYS
1	D	215	PRO
1	D	227	LYS
1	C	217	LYS
1	B	100	SER
1	B	286	VAL
1	D	100	SER
1	D	217	LYS
1	D	286	VAL
1	A	50	PRO
1	A	217	LYS
1	A	286	VAL
1	C	286	VAL
1	A	199	GLU
1	C	320	PHE
1	C	50	PRO
1	D	214	ASN
1	C	101	GLY

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	267/269 (99%)	227 (85%)	40 (15%)	4	12
1	B	267/269 (99%)	238 (89%)	29 (11%)	9	26
1	C	267/269 (99%)	233 (87%)	34 (13%)	6	18
1	D	267/269 (99%)	237 (89%)	30 (11%)	9	25
All	All	1068/1076 (99%)	935 (88%)	133 (12%)	7	19

All (133) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	15	ILE
1	A	17	GLU
1	A	19	GLU

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Mol	Chain	Res	Type
1	A	22	THR
1	A	28	VAL
1	A	31	ARG
1	A	42	LEU
1	A	52	LYS
1	A	54	LYS
1	A	57	LEU
1	A	79	VAL
1	A	83	VAL
1	A	107	GLU
1	A	116	VAL
1	A	137	ASN
1	A	151	VAL
1	A	153	THR
1	A	155	LYS
1	A	189	LEU
1	A	205	LYS
1	A	216	LEU
1	A	225	LYS
1	A	251	ILE
1	A	257	CYS
1	A	259	LEU
1	A	260	VAL
1	A	271	ILE
1	A	276	LEU
1	A	285	ILE
1	A	290	LYS
1	A	304	VAL
1	A	306	THR
1	A	308	ILE
1	A	313	LEU
1	A	318	GLU
1	A	325	LYS
1	A	331	ARG
1	A	335	THR
1	A	336	LEU
1	C	8	GLN
1	C	15	ILE
1	C	22	THR
1	C	24	SER
1	C	52	LYS
1	C	57	LEU

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Mol	Chain	Res	Type
1	C	75	THR
1	C	107	GLU
1	C	108	HIS
1	C	137	ASN
1	C	148	CYS
1	C	151	VAL
1	C	153	THR
1	C	155	LYS
1	C	178	HIS
1	C	189	LEU
1	C	205	LYS
1	C	216	LEU
1	C	222	LYS
1	C	225	LYS
1	C	244	PHE
1	C	259	LEU
1	C	260	VAL
1	C	267	MET
1	C	276	LEU
1	C	306	THR
1	C	308	ILE
1	C	313	LEU
1	C	321	ASP
1	C	324	LEU
1	C	325	LYS
1	C	331	ARG
1	C	335	THR
1	C	336	LEU
1	B	22	THR
1	B	33	LYS
1	B	74	VAL
1	B	83	VAL
1	B	100	SER
1	B	104	THR
1	B	137	ASN
1	B	151	VAL
1	B	153	THR
1	B	155	LYS
1	B	189	LEU
1	B	205	LYS
1	B	216	LEU
1	B	227	LYS

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Mol	Chain	Res	Type
1	B	241	LYS
1	B	251	ILE
1	B	259	LEU
1	B	265	GLU
1	B	269	ILE
1	B	276	LEU
1	B	285	ILE
1	B	288	THR
1	B	304	VAL
1	B	306	THR
1	B	313	LEU
1	B	318	GLU
1	B	324	LEU
1	B	325	LYS
1	B	335	THR
1	D	27	GLU
1	D	33	LYS
1	D	42	LEU
1	D	52	LYS
1	D	67	VAL
1	D	79	VAL
1	D	83	VAL
1	D	100	SER
1	D	104	THR
1	D	115	SER
1	D	137	ASN
1	D	151	VAL
1	D	153	THR
1	D	155	LYS
1	D	189	LEU
1	D	205	LYS
1	D	225	LYS
1	D	227	LYS
1	D	244	PHE
1	D	251	ILE
1	D	259	LEU
1	D	260	VAL
1	D	267	MET
1	D	276	LEU
1	D	304	VAL
1	D	306	THR
1	D	318	GLU

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Mol	Chain	Res	Type
1	D	325	LYS
1	D	328	ILE
1	D	336	LEU

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

Mol	Chain	Res	Type
1	A	8	GLN
1	A	293	GLN
1	A	317	ASN
1	C	94	HIS
1	C	245	GLN
1	C	249	ASN
1	C	317	ASN
1	B	46	HIS
1	B	249	ASN
1	B	293	GLN
1	B	311	GLN
1	D	102	GLN
1	D	214	ASN
1	D	249	ASN
1	D	293	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 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	SO4	A	342	-	4,4,4	0.16	0	6,6,6	0.50	0
3	BMD	A	600	-	5,5,5	0.42	0	5,5,5	0.44	0
4	SO4	B	342	-	4,4,4	0.12	0	6,6,6	0.46	0
3	BMD	B	500	-	5,5,5	0.17	0	5,5,5	1.07	0
4	SO4	C	342	-	4,4,4	0.22	0	6,6,6	0.41	0
3	BMD	C	700	-	5,5,5	0.34	0	5,5,5	0.47	0
4	SO4	D	342	-	4,4,4	0.17	0	6,6,6	0.37	0
3	BMD	D	400	-	5,5,5	0.27	0	5,5,5	1.07	1 (20%)

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	SO4	A	342	-	-	0/0/0/0	0/0/0/0
3	BMD	A	600	-	-	0/3/3/3	0/0/0/0
4	SO4	B	342	-	-	0/0/0/0	0/0/0/0
3	BMD	B	500	-	-	0/3/3/3	0/0/0/0
4	SO4	C	342	-	-	0/0/0/0	0/0/0/0
3	BMD	C	700	-	-	0/3/3/3	0/0/0/0
4	SO4	D	342	-	-	0/0/0/0	0/0/0/0
3	BMD	D	400	-	-	0/3/3/3	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	400	BMD	C2-C1-CA	-2.35	108.33	113.37

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

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	337/339 (99%)	-0.07	0 100 100	15, 37, 59, 64	0
1	B	337/339 (99%)	-0.09	0 100 100	15, 37, 58, 64	0
1	C	337/339 (99%)	-0.10	0 100 100	16, 37, 59, 64	0
1	D	337/339 (99%)	-0.08	0 100 100	15, 37, 58, 64	0
All	All	1348/1356 (99%)	-0.08	0 100 100	15, 37, 58, 64	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	BMD	C	700	6/6	0.34	6.32	26,29,30,30	0
3	BMD	A	600	6/6	0.29	3.16	25,28,28,29	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	BMD	B	500	6/6	0.30	2.92	38,39,39,40	0
4	SO4	C	342	5/5	0.23	1.22	103,104,105,105	0
3	BMD	D	400	6/6	0.23	1.21	35,36,36,37	0
4	SO4	A	342	5/5	0.21	0.62	102,103,103,104	0
4	SO4	D	342	5/5	0.17	-0.81	90,91,92,92	0
2	ZN	B	341	1/1	0.12	-1.13	30,30,30,30	0
4	SO4	B	342	5/5	0.16	-1.81	86,86,87,87	0
2	ZN	B	340	1/1	0.08	-2.33	39,39,39,39	0
2	ZN	A	341	1/1	0.10	-2.46	31,31,31,31	0
2	ZN	C	341	1/1	0.08	-2.80	32,32,32,32	0
2	ZN	D	340	1/1	0.07	-2.99	43,43,43,43	0
2	ZN	C	340	1/1	0.07	-3.01	48,48,48,48	0
2	ZN	A	340	1/1	0.05	-3.21	50,50,50,50	0
2	ZN	D	341	1/1	0.09	-3.57	31,31,31,31	0

6.5 Other polymers ⓘ

There are no such residues in this entry.