



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:50 AM GMT

PDB ID : 3N5K
Title : Structure Of The (Sr)Ca²⁺-ATPase E2-AlF₄- Form
Authors : Bublitz, M.; Olesen, C.; Poulsen, H.; Morth, J.P.; Moller, J.V.; Nissen, P.
Deposited on : 2010-05-25
Resolution : 2.20 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

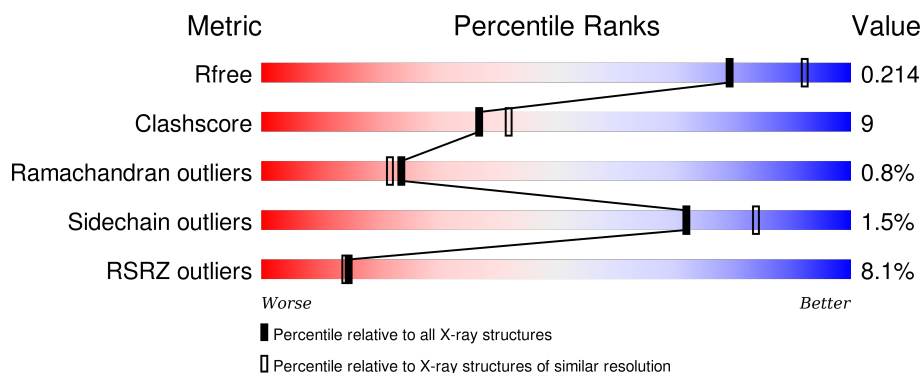
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	994	<div> <div>11%</div> <div>82%</div> <div>18%</div> </div>
1	B	994	<div> <div>5%</div> <div>82%</div> <div>17%</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
4	ALF	B	2002	-	-	X	-
6	ACT	A	3001	-	-	X	X
6	ACT	A	3002	-	-	-	X
6	ACT	B	3001	-	-	X	X
6	ACT	B	3002	-	-	X	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 16284 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.

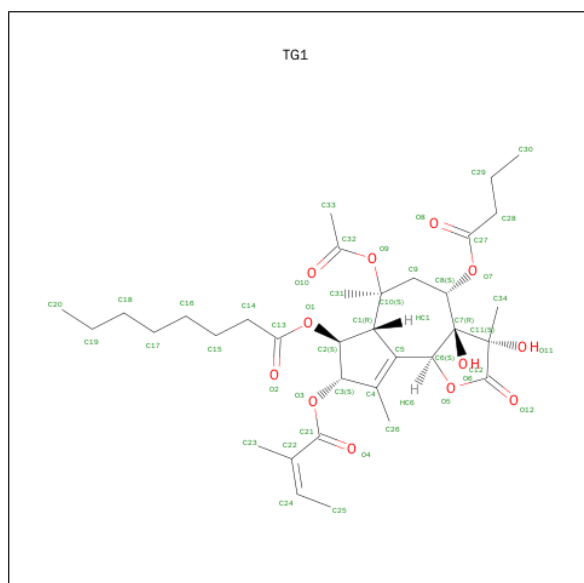
- Molecule 1 is a protein called Sarcoplasmic/endoplasmic reticulum calcium ATPase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	994	Total	C	N	O	S	0	0	0
			7671	4876	1287	1451	57			
1	B	994	Total	C	N	O	S	0	0	0
			7671	4876	1287	1451	57			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	994	GLY	ASP	SEE REMARK 999	UNP P04191
B	994	GLY	ASP	SEE REMARK 999	UNP P04191

- Molecule 2 is OCTANOIC ACID [3S-[3ALPHA, 3ABETA, 4ALPHA, 6BETA, 6ABETA, 7BETA, 8ALPHA(Z), 9BALPHA]]-6-(ACETYLOXY)-2,3,-3A,4,5,6,6A,7,8,9B-DECAHYDRO-3,3A-DIHYDROXY-3,6,9-TRIMETHYL-8-[(2-METHYL-1-OXO-2-BUTENYL)OXY]-2-OXO-4-(1-OXOBUTOXY)-AZULENO[4,5-B]FURAN-7-YL ESTER (three-letter code: TG1) (formula: C₃₄H₅₀O₁₂).

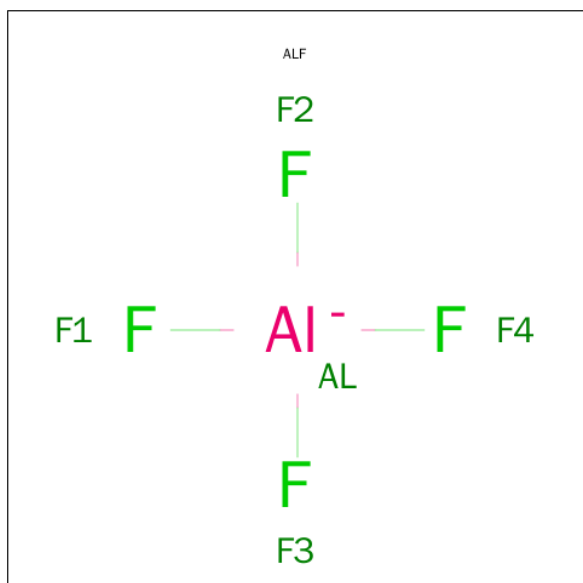


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			46	34	12		
2	B	1	Total	C	O	0	0
			46	34	12		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is TETRAFLUOROALUMINATE ION (three-letter code: ALF) (formula: AlF_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	Al	F	0	0
			5	1	4		
4	B	1	Total	Al	F	0	0
			5	1	4		

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

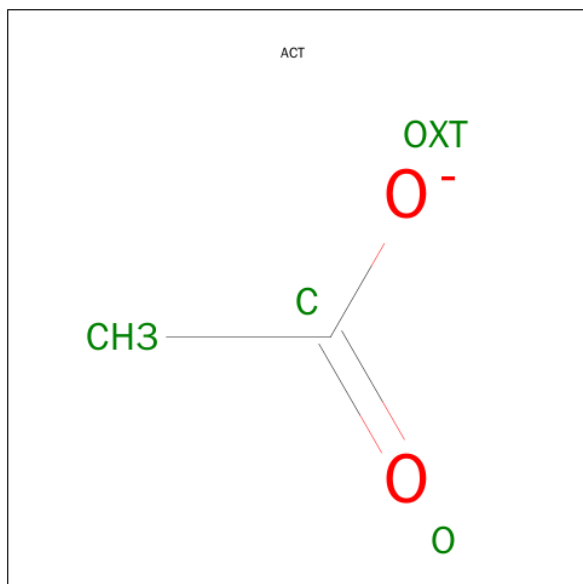
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	K	0	0
			1	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total K 1 1	0	0

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

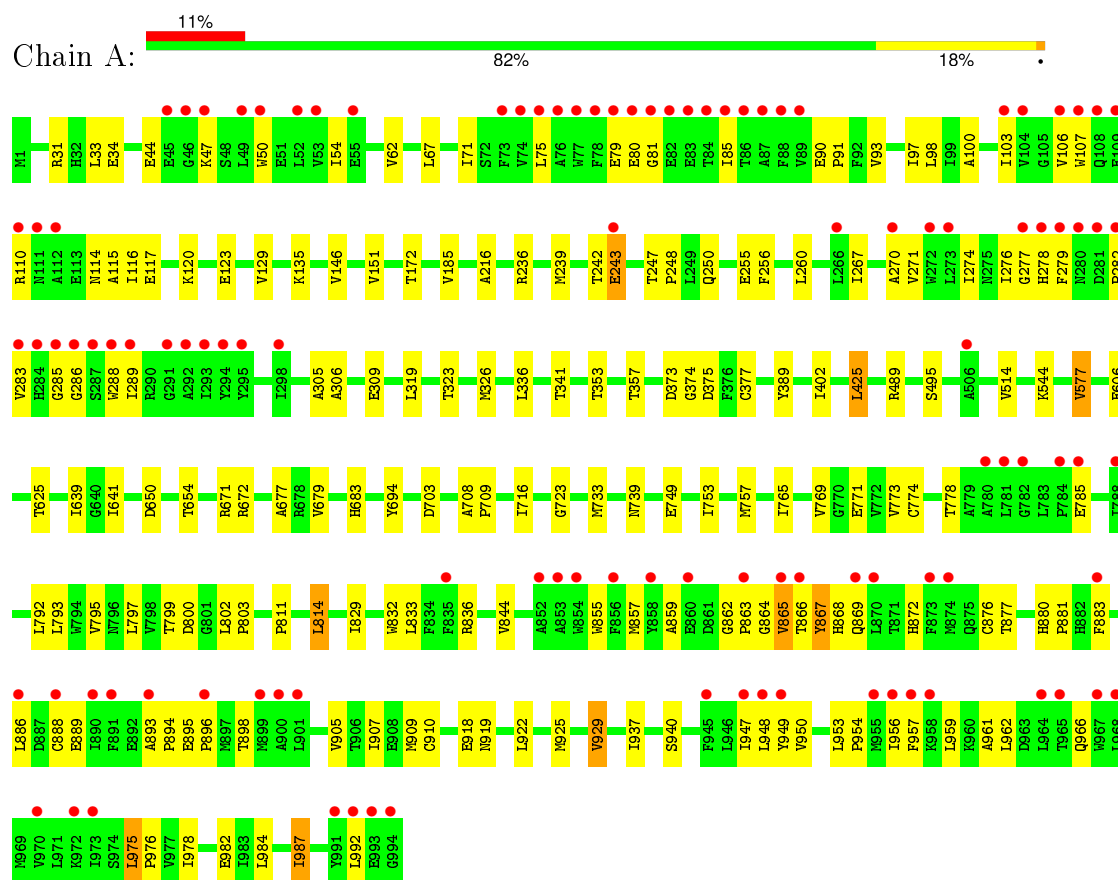
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	440	Total O 440 440	0	0
7	B	380	Total O 380 380	0	0

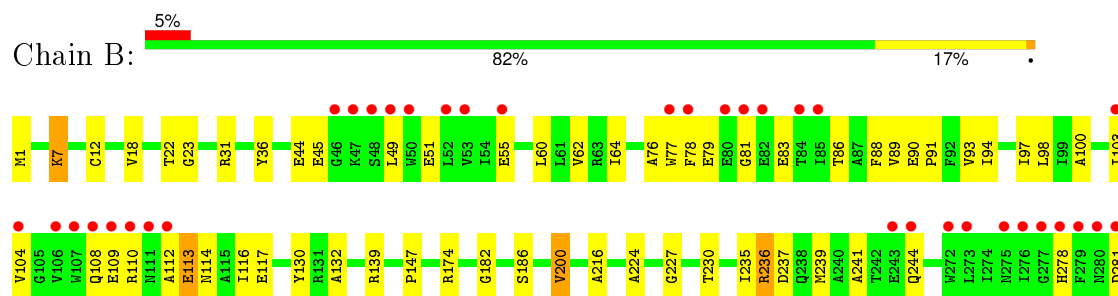
3 Residue-property plots

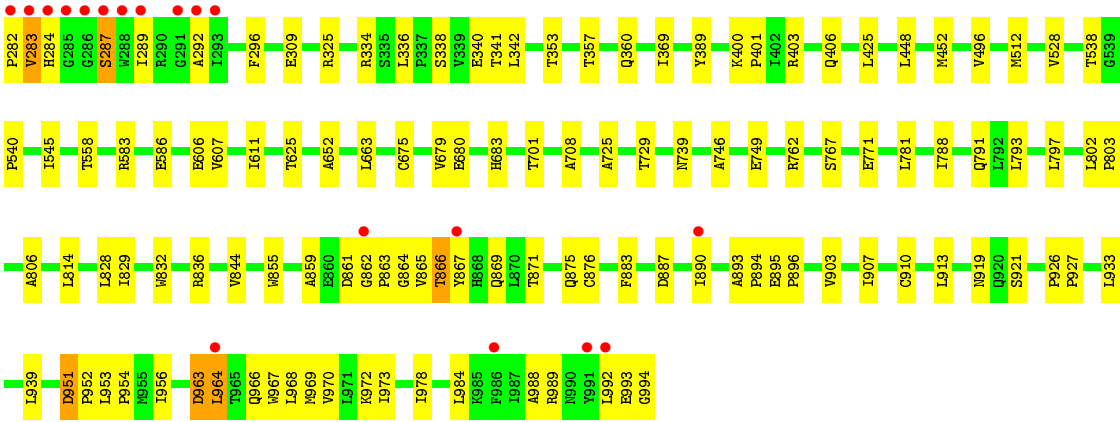
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Sarcoplasmic/endoplasmic reticulum calcium ATPase 1



- Molecule 1: Sarcoplasmic/endoplasmic reticulum calcium ATPase 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.93Å 109.42Å 276.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	72.00 – 2.20 72.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (72.00-2.20) 99.7 (72.00-2.20)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.20Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.6.1_357)	Depositor
R, R_{free}	0.184 , 0.216 0.183 , 0.214	Depositor DCC
R_{free} test set	7823 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	37.6	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 50.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 156463 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16284	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.35% 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: ALF, MG, TG1, K, ACT

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	0.51	0/7812	0.58	1/10592 (0.0%)
1	B	0.49	0/7812	0.59	1/10592 (0.0%)
All	All	0.50	0/15624	0.59	2/21184 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	7	LYS	CA-CB-CG	5.48	125.45	113.40
1	A	425	LEU	CA-CB-CG	5.27	127.42	115.30

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	7671	0	7764	138	0
1	B	7671	0	7764	146	0
2	A	46	0	50	7	0
2	B	46	0	50	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	5	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	5	0	0	2	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	8	0	6	3	0
6	B	8	0	6	5	0
7	A	440	0	0	3	0
7	B	380	0	0	0	0
All	All	16284	0	15640	293	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 9.

All (293) 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)
2:A:1000:TG1:H232	2:A:1000:TG1:H161	1.40	0.98
1:B:1:MET:HG3	1:B:224:ALA:O	1.65	0.94
1:A:242:THR:HA	1:A:243:GLU:CB	1.98	0.94
1:A:909:MET:HE3	1:A:937:ILE:HG23	1.57	0.84
1:B:762:ARG:HG2	1:B:829:ILE:HD11	1.64	0.80
1:B:1:MET:N	1:B:7:LYS:NZ	2.30	0.80
1:A:260:LEU:HD11	1:A:306:ALA:HB1	1.66	0.77
1:A:671:ARG:HD2	1:A:694:TYR:CE2	2.20	0.77
1:A:90:GLU:HB3	1:A:91:PRO:HD3	1.67	0.76
1:A:242:THR:HA	1:A:243:GLU:HB3	1.68	0.76
1:B:216:ALA:HB1	6:B:3001:ACT:H2	1.68	0.76
1:A:242:THR:HA	1:A:243:GLU:HB2	1.66	0.75
1:A:862:GLY:H	1:A:863:PRO:HD3	1.51	0.75
1:B:861:ASP:N	1:B:862:GLY:HA3	2.02	0.74
1:A:774:CYS:O	1:A:778:THR:HG22	1.86	0.74
1:B:239:MET:HE2	1:B:708:ALA:CB	2.19	0.71
1:A:855:TRP:CE3	1:A:896:PRO:HG3	2.25	0.71
1:B:90:GLU:HB3	1:B:91:PRO:HD3	1.73	0.71
1:A:267:ILE:HD11	2:A:1000:TG1:H182	1.72	0.70
1:B:338:SER:HA	1:B:341:THR:CG2	2.21	0.70
1:A:31:ARG:HH21	1:A:34:GLU:HG3	1.57	0.70
1:A:247:THR:HG23	1:A:250:GLN:H	1.56	0.69
1:A:671:ARG:HD2	1:A:694:TYR:CZ	2.27	0.69
1:B:239:MET:CE	1:B:708:ALA:HB3	2.23	0.69
1:B:7:LYS:HD2	1:B:12:CYS:SG	2.32	0.69
1:A:31:ARG:NH2	1:A:34:GLU:HG3	2.09	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ASN:HB3	1:A:117:GLU:HG2	1.73	0.68
1:B:338:SER:HA	1:B:341:THR:HG22	1.77	0.67
1:B:230:THR:OG1	6:B:3002:ACT:H2	1.94	0.67
1:B:762:ARG:HG2	1:B:829:ILE:CD1	2.24	0.67
1:B:89:VAL:O	1:B:93:VAL:HG23	1.95	0.67
1:B:1:MET:H2	1:B:7:LYS:NZ	1.92	0.66
1:A:814:LEU:H	1:A:814:LEU:HD12	1.59	0.66
1:B:963:ASP:HB3	1:B:966:GLN:H	1.62	0.64
1:B:1:MET:H1	1:B:7:LYS:NZ	1.96	0.63
1:A:606:GLU:HG3	1:A:739:ASN:OD1	1.99	0.63
1:B:863:PRO:HB2	1:B:864:GLY:HA2	1.81	0.63
1:B:887:ASP:O	1:B:890:ILE:HG12	2.00	0.62
2:A:1000:TG1:H333	2:A:1000:TG1:HC91	1.82	0.62
1:B:282:PRO:HD2	1:B:284:HIS:CE1	2.33	0.62
1:B:583:ARG:O	1:B:586:GLU:HG2	2.00	0.62
1:A:285:GLY:N	1:A:286:GLY:HA3	2.13	0.62
1:B:1:MET:N	1:B:7:LYS:HZ1	1.98	0.62
1:A:242:THR:CA	1:A:243:GLU:CB	2.77	0.62
1:A:866:THR:HB	1:A:869:GLN:HB2	1.82	0.61
1:A:905:VAL:O	1:A:909:MET:HG2	2.00	0.61
1:B:100:ALA:O	1:B:103:ILE:HG12	2.00	0.61
1:A:802:LEU:HB2	1:A:803:PRO:HD3	1.82	0.61
1:B:910:CYS:HB3	1:B:978:ILE:HG13	1.82	0.61
1:A:319:LEU:HB3	1:A:336:LEU:HD22	1.82	0.60
1:B:200:VAL:HG22	1:B:680:GLU:HG3	1.84	0.60
1:B:836:ARG:HG2	1:B:984:LEU:HB3	1.82	0.60
1:B:867:TYR:CZ	1:B:871:THR:HG21	2.37	0.59
1:A:97:ILE:HD11	1:A:797:LEU:HD22	1.83	0.59
1:B:114:ASN:HD22	1:B:117:GLU:H	1.50	0.59
1:B:876:CYS:HA	1:B:883:PHE:CD2	2.37	0.59
1:B:952:PRO:O	1:B:956:ILE:HG13	2.04	0.58
2:B:1000:TG1:C33	2:B:1000:TG1:H313	2.33	0.58
1:B:76:ALA:O	1:B:79:GLU:HB3	2.03	0.58
1:B:988:ALA:HA	1:B:992:LEU:HB2	1.84	0.58
1:A:309:GLU:HG3	1:A:793:LEU:HD23	1.85	0.58
1:A:864:GLY:O	1:A:866:THR:HG23	2.04	0.58
1:B:865:VAL:HG12	1:B:869:GLN:HB2	1.86	0.58
1:B:855:TRP:CE3	1:B:896:PRO:HG3	2.40	0.57
1:A:242:THR:CA	1:A:243:GLU:HB2	2.33	0.57
1:A:844:VAL:HG22	1:A:907:ILE:HG21	1.87	0.57
1:B:353:THR:HA	1:B:357:THR:OG1	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:866:THR:HG23	1:B:869:GLN:H	1.70	0.57
1:A:247:THR:CG2	1:A:250:GLN:H	2.17	0.57
1:B:239:MET:HE3	1:B:708:ALA:HB3	1.86	0.56
1:B:239:MET:CE	1:B:708:ALA:CB	2.80	0.56
1:B:51:GLU:O	1:B:55:GLU:HG3	2.05	0.56
1:A:893:ALA:O	1:A:896:PRO:HD2	2.05	0.56
1:A:278:HIS:O	1:A:282:PRO:HB3	2.05	0.56
1:A:863:PRO:CB	1:A:864:GLY:HA2	2.34	0.56
1:A:120:LYS:HG2	1:A:123:GLU:OE2	2.06	0.56
1:B:863:PRO:HB2	1:B:864:GLY:CA	2.36	0.56
1:A:679:VAL:HB	1:A:683:HIS:HB2	1.86	0.56
1:A:863:PRO:HB2	1:A:864:GLY:HA2	1.88	0.55
1:B:844:VAL:HG22	1:B:907:ILE:HG21	1.88	0.55
1:B:216:ALA:CB	6:B:3001:ACT:H2	2.36	0.55
1:B:963:ASP:H	1:B:966:GLN:HB2	1.70	0.55
2:A:1000:TG1:C23	2:A:1000:TG1:H161	2.27	0.55
1:B:1:MET:H2	1:B:7:LYS:CE	2.20	0.55
1:B:227:GLY:O	6:B:3002:ACT:H3	2.07	0.55
1:A:120:LYS:HA	1:A:123:GLU:OE2	2.07	0.55
1:B:893:ALA:O	1:B:896:PRO:HD2	2.07	0.55
1:A:829:ILE:HD13	2:A:1000:TG1:H333	1.89	0.54
1:A:247:THR:OG1	1:A:248:PRO:HD2	2.07	0.54
1:A:855:TRP:HA	1:A:859:ALA:HB2	1.88	0.54
1:A:880:HIS:HD2	1:A:888:CYS:SG	2.30	0.54
1:A:107:TRP:HA	1:A:110:ARG:HD3	1.89	0.54
1:B:963:ASP:HB2	1:B:966:GLN:CG	2.38	0.54
2:B:1000:TG1:HC91	2:B:1000:TG1:H333	1.89	0.53
1:B:1:MET:H1	1:B:7:LYS:HZ3	1.55	0.53
1:B:832:TRP:CD1	1:B:988:ALA:HB2	2.44	0.53
1:A:270:ALA:O	1:A:274:ILE:HG12	2.09	0.53
1:A:855:TRP:HA	1:A:859:ALA:CB	2.39	0.53
1:B:239:MET:HE2	1:B:708:ALA:HB1	1.90	0.52
1:B:139:ARG:HB2	1:B:139:ARG:NH1	2.23	0.52
1:A:80:GLU:HG3	1:A:81:GLY:H	1.75	0.52
1:B:325:ARG:NH1	1:B:749:GLU:OE2	2.38	0.52
1:A:862:GLY:N	1:A:863:PRO:HD3	2.23	0.52
1:B:887:ASP:HB3	1:B:890:ILE:HD11	1.91	0.52
1:B:953:LEU:HB2	1:B:954:PRO:HD3	1.90	0.52
1:A:947:ILE:HG22	1:A:959:LEU:HD12	1.91	0.52
1:A:984:LEU:HA	1:A:987:ILE:HD11	1.92	0.52
1:B:97:ILE:HG13	1:B:797:LEU:HD21	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:MET:SD	1:A:749:GLU:HG2	2.50	0.52
1:A:863:PRO:CD	1:A:864:GLY:HA2	2.40	0.51
1:A:949:TYR:CZ	1:A:961:ALA:HB1	2.45	0.51
1:B:1:MET:H2	1:B:7:LYS:HZ1	1.57	0.51
1:B:309:GLU:HG3	1:B:793:LEU:CD2	2.40	0.51
1:A:267:ILE:O	1:A:271:VAL:HG23	2.09	0.51
1:A:894:PRO:HD2	1:A:895:GLU:OE1	2.10	0.51
1:B:836:ARG:HA	1:B:984:LEU:HD13	1.93	0.51
1:A:832:TRP:HZ2	1:A:987:ILE:HD11	1.75	0.51
1:A:909:MET:CE	1:A:937:ILE:HA	2.41	0.51
1:A:918:GLU:HG2	1:A:919:ASN:OD1	2.11	0.51
1:A:953:LEU:HB2	1:A:954:PRO:HD3	1.92	0.51
1:B:802:LEU:HB2	1:B:803:PRO:HD3	1.93	0.51
1:A:79:GLU:HG2	1:A:80:GLU:N	2.25	0.50
1:A:79:GLU:HG2	1:A:80:GLU:H	1.77	0.50
1:B:109:GLU:O	1:B:109:GLU:HG2	2.11	0.50
1:B:969:MET:O	1:B:973:ILE:HG13	2.12	0.50
1:B:369:ILE:HG13	1:B:528:VAL:HG13	1.93	0.50
1:B:863:PRO:HB2	1:B:865:VAL:N	2.27	0.50
1:A:93:VAL:O	1:A:97:ILE:HG12	2.11	0.50
1:B:867:TYR:CE2	1:B:871:THR:HG21	2.47	0.50
1:B:200:VAL:HG22	1:B:680:GLU:CG	2.41	0.49
1:B:77:TRP:HD1	1:B:78:PHE:CE1	2.30	0.49
1:B:863:PRO:CB	1:B:864:GLY:HA2	2.42	0.49
1:B:292:ALA:O	1:B:296:PHE:HD1	1.95	0.49
1:A:172:THR:OG1	6:A:3001:ACT:H1	2.13	0.49
1:A:100:ALA:O	1:A:103:ILE:HG12	2.13	0.49
1:B:227:GLY:O	6:B:3002:ACT:CH3	2.61	0.49
1:B:287:SER:C	1:B:289:ILE:H	2.16	0.49
1:A:305:ALA:HB2	1:A:792:LEU:HD13	1.95	0.49
1:A:866:THR:O	1:A:867:TYR:HB2	2.12	0.49
1:B:448:LEU:O	1:B:452:MET:HG3	2.13	0.49
1:B:963:ASP:CB	1:B:966:GLN:H	2.26	0.48
1:A:389:TYR:HB3	1:A:425:LEU:HD11	1.96	0.48
1:A:811:PRO:HG2	1:A:929:VAL:HG13	1.94	0.48
1:B:88:PHE:O	1:B:91:PRO:HD2	2.14	0.48
1:A:489:ARG:HD3	7:A:1316:HOH:O	2.13	0.48
1:A:185:VAL:HG13	1:A:185:VAL:O	2.12	0.48
1:A:639:ILE:HD11	1:A:641:ILE:HD12	1.95	0.48
1:A:31:ARG:O	1:A:34:GLU:HG2	2.13	0.48
1:B:992:LEU:O	1:B:994:GLY:HA3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:3002:ACT:H1	7:A:1082:HOH:O	2.14	0.48
1:B:919:ASN:O	1:B:989:ARG:HD3	2.14	0.48
1:A:256:PHE:CZ	1:A:765:ILE:HD13	2.49	0.48
1:A:795:VAL:HA	1:A:799:THR:HB	1.96	0.48
2:A:1000:TG1:H191	2:A:1000:TG1:H162	1.52	0.48
1:B:113:GLU:HG2	1:B:729:THR:HG22	1.94	0.48
1:A:950:VAL:O	1:A:953:LEU:HB2	2.14	0.47
1:B:44:GLU:HG3	1:B:45:GLU:N	2.29	0.47
1:B:114:ASN:HD21	1:B:116:ILE:HB	1.79	0.47
1:B:1:MET:N	1:B:7:LYS:HZ3	2.06	0.47
1:B:90:GLU:O	1:B:94:ILE:HG13	2.14	0.47
1:A:857:MET:HA	1:A:865:VAL:HA	1.96	0.47
1:A:44:GLU:HB3	1:A:114:ASN:HD21	1.78	0.47
1:B:369:ILE:HD11	1:B:545:ILE:HD11	1.96	0.47
1:B:859:ALA:O	1:B:862:GLY:HA3	2.15	0.47
1:A:276:ILE:O	1:A:279:PHE:HB3	2.14	0.47
1:A:256:PHE:CE1	1:A:765:ILE:HD13	2.49	0.47
1:B:60:LEU:O	1:B:64:ILE:HG12	2.14	0.47
1:B:802:LEU:HD13	1:B:939:LEU:HD23	1.97	0.47
1:B:236:ARG:HG3	1:B:237:ASP:N	2.28	0.47
1:A:373:ASP:O	1:A:375:ASP:N	2.47	0.47
1:A:44:GLU:HB2	1:A:116:ILE:HD12	1.95	0.47
1:B:725:ALA:O	1:B:729:THR:HG23	2.15	0.47
1:A:962:LEU:HB3	1:A:966:GLN:HB2	1.96	0.47
1:B:62:VAL:HG13	1:B:98:LEU:HD22	1.96	0.47
1:B:88:PHE:C	1:B:91:PRO:HD2	2.34	0.47
1:A:309:GLU:HG3	1:A:793:LEU:CD2	2.45	0.47
1:B:910:CYS:O	1:B:913:LEU:HB2	2.15	0.46
1:A:577:VAL:HG11	7:A:1375:HOH:O	2.14	0.46
1:A:495:SER:HB3	1:A:514:VAL:HG22	1.96	0.46
1:A:863:PRO:N	1:A:864:GLY:HA2	2.30	0.46
1:A:864:GLY:O	1:A:865:VAL:C	2.54	0.46
1:A:898:THR:HG23	1:A:948:LEU:HD21	1.97	0.46
1:B:342:LEU:HD23	1:B:746:ALA:HB1	1.98	0.46
1:A:876:CYS:HA	1:A:883:PHE:CD2	2.50	0.46
1:A:910:CYS:HB3	1:A:978:ILE:HG13	1.98	0.46
1:A:909:MET:CE	1:A:940:SER:HB2	2.46	0.46
1:B:963:ASP:HB2	1:B:966:GLN:HG3	1.98	0.46
1:B:876:CYS:HA	1:B:883:PHE:CE2	2.51	0.46
1:B:403:ARG:HB2	1:B:406:GLN:HG2	1.97	0.46
1:A:863:PRO:N	1:A:864:GLY:CA	2.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:GLY:C	1:A:279:PHE:H	2.18	0.46
1:A:353:THR:HA	1:A:357:THR:OG1	2.16	0.46
1:B:865:VAL:HG12	1:B:866:THR:H	1.80	0.45
2:B:1000:TG1:H313	2:B:1000:TG1:H333	1.98	0.45
1:A:115:ALA:HB1	1:A:239:MET:CE	2.46	0.45
1:A:650:ASP:O	1:A:672:ARG:HD2	2.17	0.45
1:B:862:GLY:H	1:B:863:PRO:HD3	1.82	0.45
1:A:129:VAL:HG12	1:A:151:VAL:HG22	1.99	0.45
1:A:90:GLU:HB3	1:A:91:PRO:CD	2.43	0.44
1:A:62:VAL:HG13	1:A:98:LEU:HD22	1.99	0.44
1:A:50:TRP:O	1:A:54:ILE:HG12	2.18	0.44
1:B:104:VAL:O	1:B:108:GLN:HG2	2.17	0.44
1:A:377:CYS:HB3	1:A:544:LYS:HG2	2.00	0.44
1:B:993:GLU:HA	1:B:994:GLY:HA3	1.75	0.44
1:A:880:HIS:N	1:A:881:PRO:CD	2.80	0.44
1:B:538:THR:OG1	1:B:540:PRO:HD2	2.18	0.44
1:B:558:THR:O	1:B:558:THR:HG22	2.18	0.44
1:B:110:ARG:HH22	1:B:112:ALA:HB2	1.83	0.44
1:B:235:ILE:CG2	1:B:239:MET:HE1	2.46	0.44
1:B:607:VAL:O	1:B:611:ILE:HG12	2.18	0.44
1:B:894:PRO:HD2	1:B:895:GLU:OE1	2.18	0.43
1:B:282:PRO:O	1:B:283:VAL:HB	2.18	0.43
2:B:1000:TG1:H171	2:B:1000:TG1:H141	1.62	0.43
1:B:22:THR:HG22	1:B:132:ALA:HB2	1.99	0.43
1:B:968:LEU:O	1:B:972:LYS:HG2	2.18	0.43
1:A:893:ALA:HA	1:A:894:PRO:HD3	1.83	0.43
1:B:606:GLU:HG3	1:B:739:ASN:OD1	2.18	0.43
1:B:893:ALA:HA	1:B:894:PRO:HD3	1.81	0.43
1:B:283:VAL:CG2	1:B:875:GLN:HE21	2.31	0.43
1:A:975:LEU:N	1:A:976:PRO:CD	2.82	0.43
1:B:7:LYS:HB2	1:B:12:CYS:SG	2.59	0.43
1:A:880:HIS:CD2	1:A:888:CYS:SG	3.11	0.43
1:B:23:GLY:HA3	1:B:130:TYR:O	2.18	0.43
1:B:281:ASP:N	1:B:282:PRO:HD3	2.32	0.43
1:A:402:ILE:HD12	1:A:402:ILE:C	2.39	0.43
1:B:967:TRP:O	1:B:970:VAL:HB	2.17	0.43
1:B:336:LEU:HD23	1:B:336:LEU:HA	1.88	0.43
1:B:855:TRP:HA	1:B:859:ALA:CB	2.49	0.43
1:B:235:ILE:CG2	1:B:239:MET:CE	2.96	0.43
1:B:903:VAL:O	1:B:907:ILE:HG13	2.18	0.43
1:A:800:ASP:C	1:A:803:PRO:HD2	2.38	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:VAL:O	1:A:110:ARG:HG3	2.19	0.43
1:B:972:LYS:HD3	1:B:972:LYS:HA	1.78	0.43
1:A:922:LEU:HA	1:A:925:MET:O	2.19	0.43
1:B:855:TRP:HA	1:B:859:ALA:HB2	2.01	0.42
1:A:895:GLU:N	1:A:896:PRO:CD	2.82	0.42
1:A:305:ALA:HB1	1:A:771:GLU:HB3	1.99	0.42
1:A:883:PHE:HB3	1:A:886:LEU:HD11	2.01	0.42
1:B:652:ALA:HA	1:B:675:CYS:O	2.19	0.42
1:B:964:LEU:HG	1:B:964:LEU:H	1.39	0.42
1:B:836:ARG:HG2	1:B:984:LEU:HD13	2.02	0.42
1:B:182:GLY:HA3	4:B:2002:ALF:F1	2.09	0.42
1:A:85:ILE:N	1:A:85:ILE:HD12	2.33	0.42
1:A:625:THR:HA	4:A:2002:ALF:F4	2.09	0.42
1:A:33:LEU:HD13	1:A:146:VAL:CG1	2.49	0.42
1:B:389:TYR:HB3	1:B:425:LEU:HD21	2.01	0.42
1:A:31:ARG:HA	1:A:34:GLU:HG2	2.00	0.42
1:A:769:VAL:O	1:A:773:VAL:HG23	2.20	0.42
1:A:135:LYS:HD2	1:A:135:LYS:HA	1.88	0.42
1:B:83:GLU:HG2	1:B:86:THR:OG1	2.19	0.42
1:A:671:ARG:CD	1:A:694:TYR:CZ	2.99	0.42
1:A:753:ILE:O	1:A:757:MET:HG3	2.20	0.42
1:B:926:PRO:HA	1:B:927:PRO:HD3	1.94	0.42
1:A:654:THR:HA	1:A:677:ALA:O	2.20	0.42
1:A:80:GLU:HG3	1:A:81:GLY:N	2.34	0.42
1:A:71:ILE:O	1:A:75:LEU:HD13	2.20	0.41
1:B:788:ILE:HG12	1:B:791:GLN:OE1	2.20	0.41
1:A:341:THR:HG22	1:A:716:ILE:HD11	2.01	0.41
1:A:67:LEU:O	1:A:71:ILE:HG13	2.20	0.41
1:B:400:LYS:HA	1:B:401:PRO:HD3	1.92	0.41
1:B:36:TYR:CG	1:B:147:PRO:HG2	2.56	0.41
1:B:679:VAL:HB	1:B:683:HIS:HB2	2.01	0.41
1:B:1:MET:HE3	1:B:1:MET:HB3	1.40	0.41
1:B:951:ASP:OD1	1:B:952:PRO:HA	2.20	0.41
1:B:139:ARG:HH11	1:B:139:ARG:HB2	1.85	0.41
1:A:836:ARG:HG3	1:A:984:LEU:HD13	2.02	0.41
1:A:216:ALA:HB1	6:A:3001:ACT:H2	2.03	0.41
1:B:174:ARG:HB3	1:B:186:SER:HB3	2.02	0.41
1:A:708:ALA:HB3	1:A:709:PRO:HD3	2.02	0.41
1:A:868:HIS:CE1	1:A:872:HIS:CE1	3.09	0.41
1:A:288:TRP:CD1	1:A:289:ILE:HG13	2.56	0.41
1:A:898:THR:CG2	1:A:948:LEU:HD21	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:814:LEU:N	1:B:814:LEU:HD12	2.36	0.41
1:A:984:LEU:HA	1:A:987:ILE:CD1	2.50	0.41
1:A:833:LEU:O	1:A:836:ARG:HB3	2.21	0.41
1:A:909:MET:HE1	1:A:937:ILE:HA	2.01	0.41
1:B:309:GLU:HG3	1:B:793:LEU:HD23	2.02	0.41
1:B:94:ILE:HG12	1:B:793:LEU:HD11	2.03	0.41
1:B:239:MET:HE2	1:B:708:ALA:HB3	1.90	0.41
1:B:496:VAL:O	1:B:512:MET:HA	2.22	0.41
1:A:956:ILE:HD11	1:A:957:PHE:CE1	2.56	0.40
1:B:767:SER:O	1:B:771:GLU:HG3	2.21	0.40
1:B:951:ASP:CG	1:B:952:PRO:HA	2.41	0.40
1:B:625:THR:HA	4:B:2002:ALF:F4	2.11	0.40
1:A:256:PHE:HB2	2:A:1000:TG1:H291	2.02	0.40
1:A:855:TRP:CZ3	1:A:896:PRO:HG3	2.56	0.40
1:A:733:MET:HB2	1:A:733:MET:HE2	1.79	0.40
1:B:278:HIS:O	1:B:278:HIS:CG	2.73	0.40
1:B:887:ASP:HB3	1:B:890:ILE:CD1	2.52	0.40
1:B:806:ALA:HB1	1:B:933:LEU:HA	2.02	0.40
1:A:947:ILE:HD13	1:A:957:PHE:CE1	2.56	0.40
1:B:828:LEU:N	1:B:828:LEU:HD12	2.36	0.40
1:A:703:ASP:HB2	1:A:723:GLY:HA3	2.04	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	992/994 (100%)	941 (95%)	41 (4%)	10 (1%)	19	16
1	B	992/994 (100%)	951 (96%)	36 (4%)	5 (0%)	34	35
All	All	1984/1988 (100%)	1892 (95%)	77 (4%)	15 (1%)	24	22

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	GLU
1	A	889	GLU
1	B	283	VAL
1	A	374	GLY
1	A	865	VAL
1	A	992	LEU
1	B	241	ALA
1	A	283	VAL
1	A	785	GLU
1	A	47	LYS
1	A	877	THR
1	B	244	GLN
1	B	287	SER
1	A	867	TYR
1	B	81	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	840/840 (100%)	831 (99%)	9 (1%)	80	89
1	B	840/840 (100%)	823 (98%)	17 (2%)	63	76
All	All	1680/1680 (100%)	1654 (98%)	26 (2%)	72	84

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	236	ARG
1	A	255	GLU
1	A	323	THR
1	A	577	VAL
1	A	814	LEU
1	A	929	VAL
1	A	975	LEU
1	A	982	GLU

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Mol	Chain	Res	Type
1	A	987	ILE
1	B	18	VAL
1	B	31	ARG
1	B	49	LEU
1	B	113	GLU
1	B	200	VAL
1	B	236	ARG
1	B	334	ARG
1	B	340	GLU
1	B	360	GLN
1	B	663	LEU
1	B	701	THR
1	B	781	LEU
1	B	866	THR
1	B	921	SER
1	B	951	ASP
1	B	963	ASP
1	B	964	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	868	HIS
1	A	880	HIS
1	B	114	ASN
1	B	284	HIS
1	B	875	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 12 ligands modelled in this entry, 4 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	$\# Z > 2$
2	TG1	A	1000	-	43,48,48	2.33	8 (18%)	42,72,72	3.97	14 (33%)
4	ALF	A	2002	-	0,4,4	0.00	-	0,6,6	0.00	-
6	ACT	A	3001	-	1,3,3	0.18	0	0,3,3	0.00	-
6	ACT	A	3002	-	1,3,3	2.56	1 (100%)	0,3,3	0.00	-
2	TG1	B	1000	-	43,48,48	2.34	8 (18%)	42,72,72	3.88	15 (35%)
4	ALF	B	2002	-	0,4,4	0.00	-	0,6,6	0.00	-
6	ACT	B	3001	-	1,3,3	0.35	0	0,3,3	0.00	-
6	ACT	B	3002	-	1,3,3	0.25	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
2	TG1	A	1000	-	-	2/33/99/99	0/3/3/3
4	ALF	A	2002	-	-	0/0/0/0	0/0/0/0
6	ACT	A	3001	-	-	0/0/0/0	0/0/0/0
6	ACT	A	3002	-	-	0/0/0/0	0/0/0/0
2	TG1	B	1000	-	-	2/33/99/99	0/3/3/3
4	ALF	B	2002	-	-	0/0/0/0	0/0/0/0
6	ACT	B	3001	-	-	0/0/0/0	0/0/0/0
6	ACT	B	3002	-	-	0/0/0/0	0/0/0/0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1000	TG1	O9-C10	-3.05	1.43	1.48
2	B	1000	TG1	O9-C10	-3.03	1.43	1.48
2	A	1000	TG1	O7-C8	-2.86	1.41	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1000	TG1	O7-C8	-2.78	1.41	1.46
2	B	1000	TG1	O3-C21	2.05	1.39	1.34
2	A	1000	TG1	C9-C8	2.20	1.54	1.52
2	B	1000	TG1	C11-C7	2.30	1.58	1.55
6	A	3002	ACT	CH3-C	2.56	1.52	1.48
2	A	1000	TG1	C11-C7	2.72	1.58	1.55
2	A	1000	TG1	C24-C22	3.36	1.53	1.30
2	B	1000	TG1	C24-C22	3.41	1.54	1.30
2	B	1000	TG1	O5-C12	3.45	1.41	1.35
2	A	1000	TG1	O5-C12	3.57	1.41	1.35
2	A	1000	TG1	O12-C12	7.16	1.38	1.20
2	B	1000	TG1	O12-C12	7.48	1.38	1.20
2	A	1000	TG1	C4-C5	9.76	1.54	1.33
2	B	1000	TG1	C4-C5	9.87	1.55	1.33

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1000	TG1	O12-C12-C11	-19.12	110.66	128.26
2	B	1000	TG1	O12-C12-C11	-17.73	111.95	128.26
2	B	1000	TG1	O5-C12-O12	-10.81	105.88	121.62
2	A	1000	TG1	O5-C12-O12	-10.80	105.89	121.62
2	B	1000	TG1	C26-C4-C3	-5.18	114.79	121.51
2	B	1000	TG1	O5-C12-C11	-4.51	103.45	110.14
2	A	1000	TG1	C23-C22-C24	-4.48	106.86	123.55
2	B	1000	TG1	C23-C22-C24	-4.15	108.11	123.55
2	B	1000	TG1	O3-C21-O4	-3.93	115.46	123.30
2	A	1000	TG1	O5-C12-C11	-3.86	104.41	110.14
2	A	1000	TG1	C7-C11-C12	-3.75	93.80	100.84
2	B	1000	TG1	C7-C11-C12	-3.68	93.93	100.84
2	A	1000	TG1	C26-C4-C3	-3.67	116.75	121.51
2	A	1000	TG1	O3-C21-O4	-3.64	116.03	123.30
2	A	1000	TG1	C23-C22-C21	-3.05	108.12	116.04
2	B	1000	TG1	C23-C22-C21	-2.67	109.11	116.04
2	B	1000	TG1	C26-C4-C5	-2.47	123.51	129.97
2	A	1000	TG1	C26-C4-C5	-2.45	123.54	129.97
2	B	1000	TG1	C3-O3-C21	-2.28	113.02	116.30
2	A	1000	TG1	O5-C6-C7	-2.27	102.07	104.08
2	A	1000	TG1	C25-C24-C22	-2.09	116.99	127.86
2	B	1000	TG1	O1-C13-O2	-2.06	118.14	123.67
2	B	1000	TG1	C25-C24-C22	-2.02	117.35	127.86
2	A	1000	TG1	O11-C11-C12	2.09	112.51	106.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1000	TG1	O7-C8-C9	2.18	111.01	107.03
2	B	1000	TG1	O11-C11-C12	2.20	112.85	106.08
2	B	1000	TG1	O7-C8-C9	2.20	111.05	107.03
2	A	1000	TG1	C10-O9-C32	6.59	135.42	121.90
2	B	1000	TG1	C10-O9-C32	6.92	136.10	121.90

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1000	TG1	C10-O9-C32-C33
2	B	1000	TG1	C10-O9-C32-C33
2	A	1000	TG1	C10-O9-C32-O10
2	B	1000	TG1	C10-O9-C32-O10

There are no ring outliers.

8 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1000	TG1	7	0
4	A	2002	ALF	1	0
6	A	3001	ACT	2	0
6	A	3002	ACT	1	0
2	B	1000	TG1	4	0
4	B	2002	ALF	2	0
6	B	3001	ACT	2	0
6	B	3002	ACT	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	994/994 (100%)	0.45	108 (10%) 7 7	19, 45, 149, 275	0
1	B	994/994 (100%)	0.07	53 (5%) 30 29	22, 49, 119, 221	0
All	All	1988/1988 (100%)	0.26	161 (8%) 15 14	19, 48, 137, 275	0

All (161) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	285	GLY	13.8
1	A	80	GLU	11.9
1	B	49	LEU	11.6
1	A	84	THR	10.4
1	B	286	GLY	9.1
1	B	80	GLU	8.9
1	A	78	PHE	8.9
1	B	81	GLY	8.7
1	A	81	GLY	8.5
1	B	112	ALA	8.5
1	A	77	TRP	7.6
1	A	110	ARG	7.5
1	B	109	GLU	7.3
1	B	283	VAL	7.3
1	A	85	ILE	7.1
1	A	76	ALA	7.0
1	A	292	ALA	6.9
1	A	283	VAL	6.6
1	A	891	PHE	6.6
1	A	111	ASN	6.5
1	A	888	CYS	6.5
1	A	992	LEU	6.4
1	B	77	TRP	6.3
1	B	280	ASN	6.2

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Mol	Chain	Res	Type	RSRZ
1	A	47	LYS	6.2
1	B	285	GLY	6.1
1	B	108	GLN	6.1
1	B	282	PRO	6.0
1	B	287	SER	5.9
1	A	286	GLY	5.9
1	A	79	GLU	5.9
1	A	82	GLU	5.8
1	B	281	ASP	5.7
1	B	78	PHE	5.7
1	B	111	ASN	5.7
1	A	50	TRP	5.6
1	A	949	TYR	5.5
1	B	289	ILE	5.5
1	B	278	HIS	5.4
1	A	883	PHE	5.2
1	A	948	LEU	5.2
1	B	288	TRP	5.0
1	A	266	LEU	5.0
1	B	84	THR	5.0
1	B	50	TRP	4.9
1	A	994	GLY	4.9
1	B	48	SER	4.9
1	B	47	LYS	4.9
1	B	276	ILE	4.8
1	B	244	GLN	4.8
1	A	279	PHE	4.7
1	A	280	ASN	4.6
1	A	853	ALA	4.6
1	A	890	ILE	4.6
1	A	781	LEU	4.6
1	A	74	VAL	4.6
1	A	106	VAL	4.5
1	A	970	VAL	4.5
1	B	275	ASN	4.4
1	B	273	LEU	4.3
1	A	873	PHE	4.3
1	A	947	ILE	4.3
1	B	279	PHE	4.3
1	A	281	ASP	4.2
1	A	835	PHE	4.2
1	A	52	LEU	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	85	ILE	4.2
1	B	110	ARG	4.1
1	A	865	VAL	4.1
1	B	107	TRP	4.1
1	B	52	LEU	4.0
1	B	46	GLY	4.0
1	A	86	THR	4.0
1	A	295	TYR	4.0
1	A	107	TRP	3.9
1	B	243	GLU	3.9
1	A	782	GLY	3.9
1	B	277	GLY	3.9
1	A	104	VAL	3.9
1	A	278	HIS	3.8
1	A	46	GLY	3.7
1	A	993	GLU	3.6
1	A	273	LEU	3.6
1	A	874	MET	3.6
1	A	88	PHE	3.6
1	A	294	TYR	3.5
1	A	856	PHE	3.5
1	A	243	GLU	3.5
1	A	288	TRP	3.5
1	B	82	GLU	3.5
1	A	854	TRP	3.5
1	A	852	ALA	3.5
1	A	293	ILE	3.5
1	A	289	ILE	3.4
1	B	106	VAL	3.4
1	A	291	GLY	3.3
1	A	957	PHE	3.3
1	A	896	PRO	3.2
1	A	298	ILE	3.2
1	A	956	ILE	3.2
1	A	858	TYR	3.2
1	A	784	PRO	3.2
1	A	45	GLU	3.1
1	A	886	LEU	3.1
1	A	108	GLN	3.1
1	A	103	ILE	3.1
1	A	991	TYR	3.1
1	A	788	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	109	GLU	3.1
1	A	83	GLU	3.1
1	B	272	TRP	3.1
1	A	284	HIS	3.0
1	A	287	SER	3.0
1	A	866	THR	3.0
1	A	282	PRO	3.0
1	A	49	LEU	2.9
1	A	900	ALA	2.9
1	A	967	TRP	2.9
1	B	293	ILE	2.9
1	A	958	LYS	2.9
1	A	89	VAL	2.9
1	A	965	THR	2.9
1	A	972	LYS	2.8
1	B	867	TYR	2.8
1	B	284	HIS	2.8
1	A	870	LEU	2.8
1	A	973	ILE	2.8
1	B	986	PHE	2.8
1	A	964	LEU	2.8
1	A	277	GLY	2.6
1	B	292	ALA	2.6
1	B	55	GLU	2.6
1	A	893	ALA	2.6
1	A	112	ALA	2.5
1	A	75	LEU	2.5
1	B	103	ILE	2.5
1	A	272	TRP	2.5
1	A	901	LEU	2.5
1	B	991	TYR	2.5
1	A	955	MET	2.5
1	A	53	VAL	2.5
1	A	860	GLU	2.4
1	B	53	VAL	2.4
1	A	506	ALA	2.4
1	A	899	MET	2.4
1	A	968	LEU	2.4
1	A	863	PRO	2.3
1	A	780	ALA	2.3
1	A	270	ALA	2.3
1	A	73	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	992	LEU	2.2
1	B	104	VAL	2.2
1	A	785	GLU	2.2
1	A	87	ALA	2.2
1	B	964	LEU	2.2
1	B	862	GLY	2.1
1	B	890	ILE	2.1
1	A	945	PHE	2.1
1	B	291	GLY	2.1
1	A	869	GLN	2.1
1	A	55	GLU	2.0

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

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

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(Å ²)	Q<0.9
6	ACT	A	3002	4/4	0.89	0.18	4.88	29,38,42,46	0
6	ACT	A	3001	4/4	0.90	0.19	4.48	35,44,45,51	0
6	ACT	B	3002	4/4	0.93	0.16	4.22	31,44,45,50	0
6	ACT	B	3001	4/4	0.92	0.16	4.19	41,48,49,55	0
3	MG	A	2001	1/1	0.98	0.16	1.78	26,26,26,26	0
3	MG	B	2001	1/1	0.99	0.14	1.67	28,28,28,28	0
4	ALF	B	2002	5/5	0.99	0.14	1.25	23,23,25,28	0
2	TG1	A	1000	46/46	0.93	0.20	0.99	52,69,119,119	0
2	TG1	B	1000	46/46	0.94	0.17	0.92	49,63,87,88	0
4	ALF	A	2002	5/5	0.99	0.14	0.60	23,23,25,27	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	K	A	2003	1/1	0.98	0.11	-0.39	36,36,36,36	0
5	K	B	2003	1/1	0.99	0.11	-0.81	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.