



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:11 AM UTC

PDB ID : 6I60 / pdb_00006i60
Title : Structure of alpha-L-rhamnosidase from Dictyoglumus thermophilum
Authors : Lafite, P.; Daniellou, R.
Deposited on : 2018-11-15
Resolution : 2.74 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

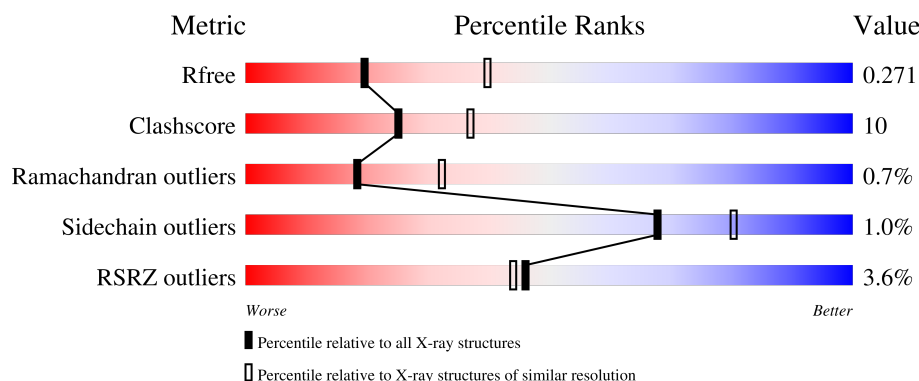
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

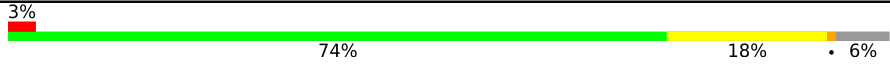
The reported resolution of this entry is 2.74 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	944	
1	B	944	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PGE	A	1001	-	-	X	-
2	PGE	A	1007	-	-	X	-
2	PGE	A	1010	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14789 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 Alpha-rhamnosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	883	Total	C	N	O	S	0	0	0
			7237	4694	1169	1356	18			
1	B	886	Total	C	N	O	S	0	0	0
			7251	4705	1174	1354	18			

There are 46 discrepancies between the modelled and reference sequences:

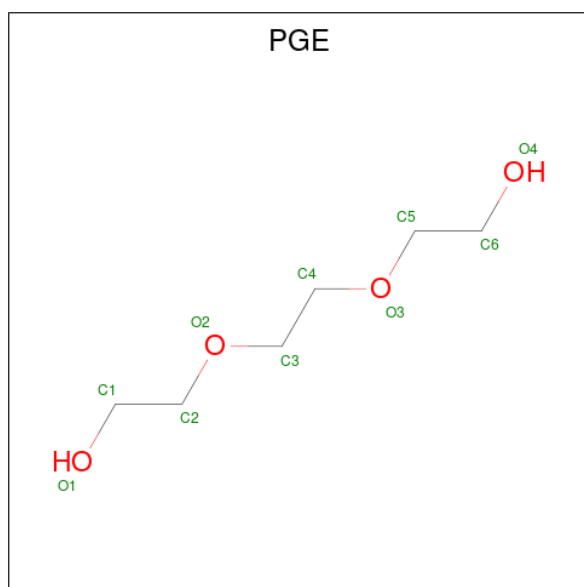
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP B5YC64
A	2	GLY	-	expression tag	UNP B5YC64
A	3	SER	-	expression tag	UNP B5YC64
A	4	SER	-	expression tag	UNP B5YC64
A	5	HIS	-	expression tag	UNP B5YC64
A	6	HIS	-	expression tag	UNP B5YC64
A	7	HIS	-	expression tag	UNP B5YC64
A	8	HIS	-	expression tag	UNP B5YC64
A	9	HIS	-	expression tag	UNP B5YC64
A	10	HIS	-	expression tag	UNP B5YC64
A	11	SER	-	expression tag	UNP B5YC64
A	12	SER	-	expression tag	UNP B5YC64
A	13	GLY	-	expression tag	UNP B5YC64
A	14	LEU	-	expression tag	UNP B5YC64
A	15	VAL	-	expression tag	UNP B5YC64
A	16	PRO	-	expression tag	UNP B5YC64
A	17	ARG	-	expression tag	UNP B5YC64
A	18	GLY	-	expression tag	UNP B5YC64
A	19	SER	-	expression tag	UNP B5YC64
A	20	HIS	-	expression tag	UNP B5YC64
A	21	MET	-	expression tag	UNP B5YC64
A	22	ALA	-	expression tag	UNP B5YC64
A	23	SER	-	expression tag	UNP B5YC64
B	1	MET	-	initiating methionine	UNP B5YC64
B	2	GLY	-	expression tag	UNP B5YC64

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Chain	Residue	Modelled	Actual	Comment	Reference
B	3	SER	-	expression tag	UNP B5YC64
B	4	SER	-	expression tag	UNP B5YC64
B	5	HIS	-	expression tag	UNP B5YC64
B	6	HIS	-	expression tag	UNP B5YC64
B	7	HIS	-	expression tag	UNP B5YC64
B	8	HIS	-	expression tag	UNP B5YC64
B	9	HIS	-	expression tag	UNP B5YC64
B	10	HIS	-	expression tag	UNP B5YC64
B	11	SER	-	expression tag	UNP B5YC64
B	12	SER	-	expression tag	UNP B5YC64
B	13	GLY	-	expression tag	UNP B5YC64
B	14	LEU	-	expression tag	UNP B5YC64
B	15	VAL	-	expression tag	UNP B5YC64
B	16	PRO	-	expression tag	UNP B5YC64
B	17	ARG	-	expression tag	UNP B5YC64
B	18	GLY	-	expression tag	UNP B5YC64
B	19	SER	-	expression tag	UNP B5YC64
B	20	HIS	-	expression tag	UNP B5YC64
B	21	MET	-	expression tag	UNP B5YC64
B	22	ALA	-	expression tag	UNP B5YC64
B	23	SER	-	expression tag	UNP B5YC64

- Molecule 2 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



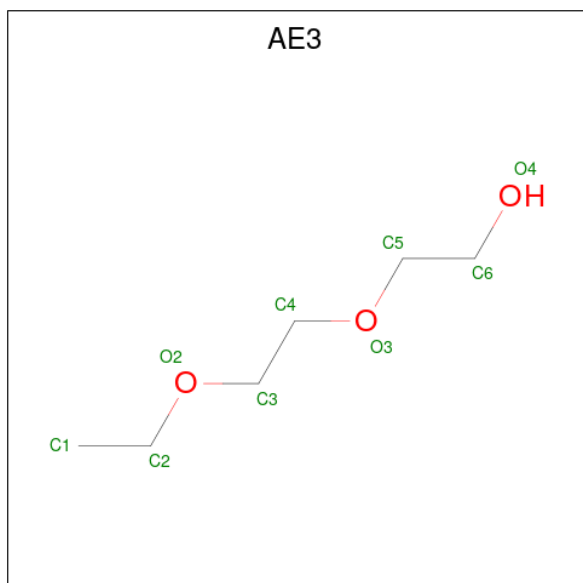
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	6	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	6	4		
2	A	1	Total	C	O	0	0
			10	6	4		
2	A	1	Total	C	O	0	0
			10	6	4		
2	A	1	Total	C	O	0	0
			10	6	4		
2	A	1	Total	C	O	0	0
			10	6	4		
2	B	1	Total	C	O	0	0
			10	6	4		
2	B	1	Total	C	O	0	0
			10	6	4		
2	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 3 is 2-(2-ETHOXYETHOXY)ETHANOL (CCD ID: AE3) (formula: C₆H₁₄O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			9	6	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			8	5	3		
3	B	1	Total	C	O	0	0
			9	6	3		
3	B	1	Total	C	O	0	0
			8	5	3		
3	B	1	Total	C	O	0	0
			6	4	2		

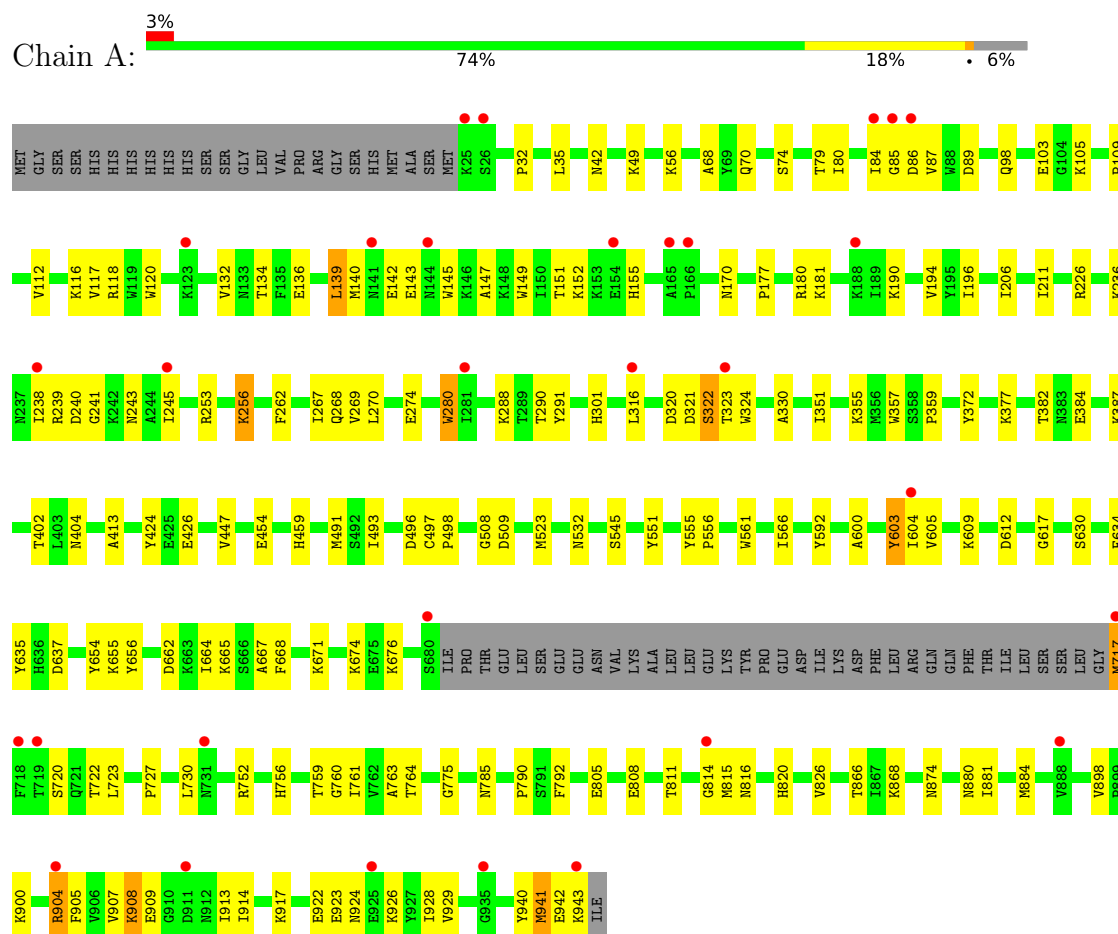
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	81	Total	O	0	0
			81	81		
4	B	70	Total	O	0	0
			70	70		

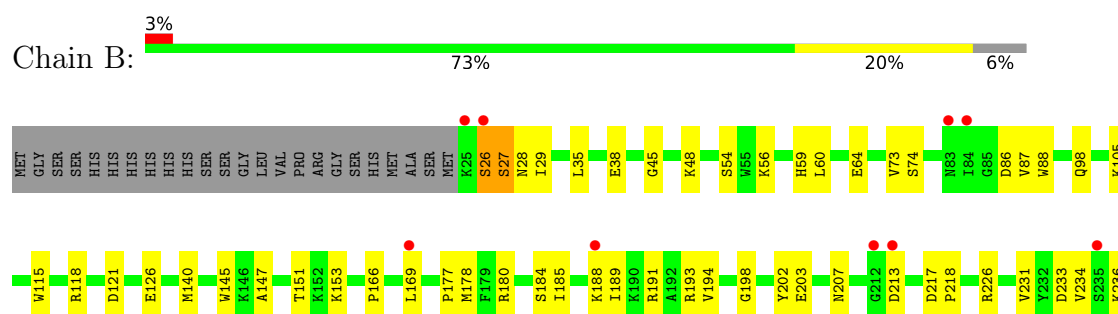
3 Residue-property plots [i](#)

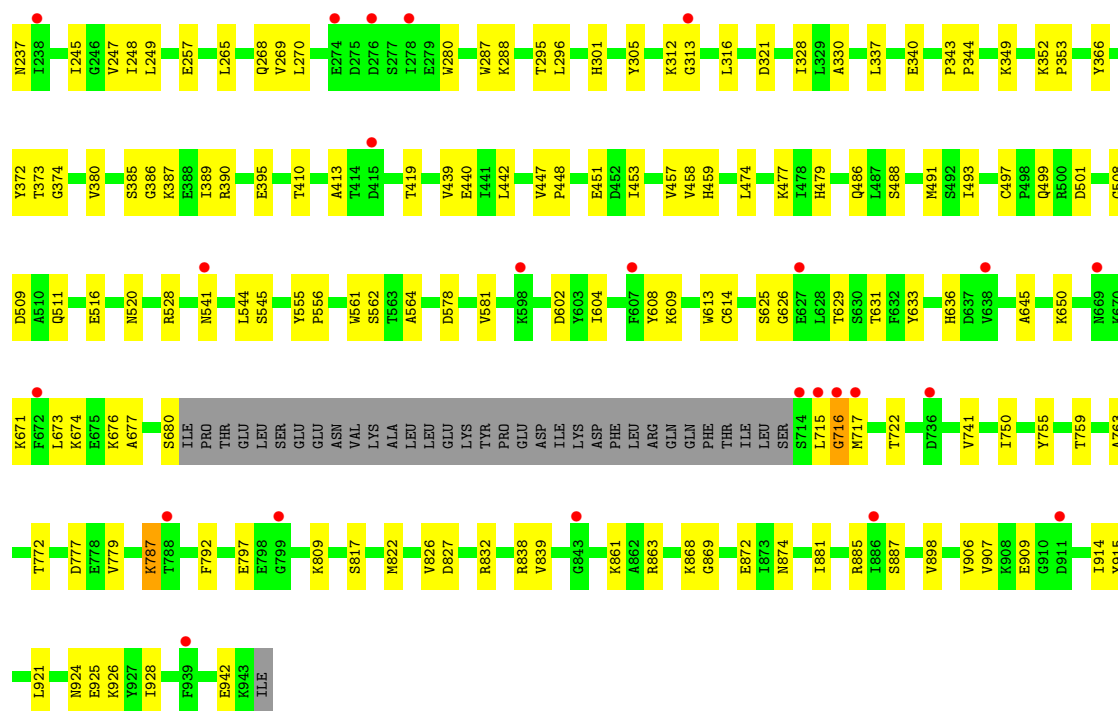
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Alpha-rhamnosidase



• Molecule 1: Alpha-rhamnosidase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	173.22Å 161.53Å 124.38Å 90.00° 132.14° 90.00°	Depositor
Resolution (Å)	43.28 – 2.74 43.28 – 2.74	Depositor EDS
% Data completeness (in resolution range)	96.9 (43.28-2.74) 97.0 (43.28-2.74)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.94 (at 2.73Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.226 , 0.270 0.230 , 0.271	Depositor DCC
R_{free} test set	3265 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	60.8	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.042 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	14789	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: PGE, AE3

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	0.33	2/7442 (0.0%)	0.34	1/10098 (0.0%)
1	B	0.18	1/7456 (0.0%)	0.31	3/10115 (0.0%)
All	All	0.27	3/14898 (0.0%)	0.33	4/20213 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	305	TYR	C-O	-6.60	1.16	1.24
1	A	656	TYR	C-O	-5.90	1.17	1.24
1	A	654	TYR	C-O	-5.21	1.18	1.24

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	26	SER	N-CA-C	7.47	117.82	108.19
1	B	27	SER	N-CA-C	6.79	121.58	113.16
1	A	603	TYR	N-CA-C	5.97	122.80	113.72
1	B	28	ASN	N-CA-C	-5.18	100.07	108.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7237	0	7030	141	0
1	B	7251	0	7059	139	0
2	A	80	0	112	30	0
2	B	30	0	42	6	0
3	A	17	0	23	1	0
3	B	23	0	32	4	0
4	A	81	0	0	2	0
4	B	70	0	0	3	0
All	All	14789	0	14298	281	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 10.

All (281) 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:256:LYS:CE	2:A:1007:PGE:H42	1.79	1.11
1:A:355:LYS:HE2	2:A:1001:PGE:C6	1.81	1.10
1:A:256:LYS:HE3	2:A:1007:PGE:C4	1.83	1.08
1:B:787:LYS:HE3	3:B:1001:AE3:H1C2	1.40	1.04
1:B:27:SER:O	1:B:29:ILE:HD12	1.55	1.04
1:A:355:LYS:HE2	2:A:1001:PGE:H6	1.41	1.00
1:A:256:LYS:HE3	2:A:1007:PGE:H42	1.01	0.99
1:A:355:LYS:NZ	2:A:1001:PGE:H62	1.76	0.99
1:A:717:MET:HG2	1:A:752:ARG:NH2	1.77	0.98
1:A:355:LYS:CE	2:A:1001:PGE:H62	1.94	0.96
1:A:256:LYS:HE2	2:A:1007:PGE:H1	1.48	0.92
1:A:355:LYS:HE2	2:A:1001:PGE:H62	1.51	0.88
1:A:717:MET:HG2	1:A:752:ARG:HH22	1.39	0.87
1:A:355:LYS:CE	2:A:1001:PGE:C6	2.54	0.85
1:A:900:LYS:HA	1:A:941:MET:HE1	1.61	0.82
1:B:27:SER:O	1:B:29:ILE:CD1	2.29	0.80
1:A:206:ILE:HG12	1:A:245:ILE:HG22	1.67	0.77
1:A:268:GLN:NE2	1:A:280:TRP:CZ3	2.53	0.77
1:A:775:GLY:HA3	2:A:1010:PGE:H2	1.68	0.76
1:A:270:LEU:HD23	1:A:280:TRP:CE3	2.22	0.75
1:A:926:LYS:HE3	2:A:1010:PGE:H22	1.66	0.75
1:A:196:ILE:HB	1:A:267:ILE:HD13	1.69	0.75
1:B:906:VAL:HG12	1:B:942:GLU:HB2	1.67	0.74
1:B:29:ILE:HD13	1:B:60:LEU:HD13	1.69	0.74
1:A:85:GLY:HA3	1:A:89:ASP:HB2	1.70	0.74
1:B:477:LYS:HZ2	2:B:1004:PGE:H52	1.53	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:904:ARG:HH12	1:A:943:LYS:CB	2.02	0.71
1:B:180:ARG:NH1	1:B:321:ASP:OD1	2.23	0.71
1:A:355:LYS:HZ3	2:A:1001:PGE:H62	1.54	0.70
1:B:194:VAL:HG22	1:B:269:VAL:HG22	1.71	0.70
1:A:181:LYS:HB3	1:A:245:ILE:HD11	1.74	0.69
1:A:926:LYS:HD2	2:A:1010:PGE:H42	1.75	0.68
1:A:35:LEU:HD21	1:A:117:VAL:HG12	1.75	0.67
1:B:410:THR:HB	1:B:501:ASP:OD2	1.96	0.66
1:A:603:TYR:C	1:A:604:ILE:HD12	2.21	0.66
1:B:366:TYR:HB2	1:B:439:VAL:HG12	1.76	0.65
1:B:184:SER:O	1:B:185:ILE:HD12	1.97	0.65
1:A:256:LYS:HE2	2:A:1007:PGE:C1	2.28	0.64
1:A:70:GLN:OE1	1:A:118:ARG:NH2	2.31	0.63
1:B:184:SER:C	1:B:185:ILE:HD12	2.23	0.63
1:B:907:VAL:HB	1:B:915:TYR:HB3	1.81	0.63
1:A:194:VAL:HG13	1:A:267:ILE:HD11	1.80	0.62
1:B:145:TRP:O	1:B:147:ALA:N	2.31	0.62
1:A:79:THR:HG23	1:A:84:ILE:HD11	1.82	0.62
1:A:238:ILE:O	1:A:243:ASN:ND2	2.31	0.62
1:A:775:GLY:HA2	2:A:1010:PGE:H6	1.81	0.62
1:A:256:LYS:CE	2:A:1007:PGE:C4	2.58	0.62
1:A:270:LEU:HA	1:A:280:TRP:HB3	1.81	0.62
1:A:604:ILE:HD12	1:A:604:ILE:N	2.15	0.62
1:B:295:THR:OG1	1:B:296:LEU:N	2.33	0.61
1:B:193:ARG:NH1	1:B:233:ASP:OD2	2.34	0.61
1:B:477:LYS:NZ	2:B:1004:PGE:H52	2.15	0.61
1:A:180:ARG:NH2	1:A:316:LEU:O	2.34	0.61
1:B:440:GLU:OE2	1:B:442:LEU:HD11	2.01	0.61
1:B:98:GLN:OE1	1:B:868:LYS:NZ	2.33	0.60
1:A:142:GLU:HA	1:A:145:TRP:CD1	2.36	0.60
1:A:322:SER:O	1:A:324:TRP:N	2.34	0.60
1:B:715:LEU:HD22	1:B:716:GLY:H	1.66	0.60
1:B:312:LYS:HD3	1:B:313:GLY:N	2.17	0.59
1:A:151:THR:OG1	1:A:330:ALA:HB2	2.02	0.59
1:B:226:ARG:NH1	1:B:340:GLU:OE1	2.37	0.58
1:A:664:ILE:O	1:A:665:LYS:C	2.46	0.58
1:A:909:GLU:HB2	1:A:914:ILE:HD13	1.84	0.58
1:B:27:SER:C	1:B:29:ILE:HD12	2.26	0.58
1:B:477:LYS:NZ	2:B:1004:PGE:O1	2.34	0.58
1:A:600:ALA:HB2	1:A:605:VAL:HG13	1.85	0.57
1:A:556:PRO:O	1:A:609:LYS:NZ	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:528:ARG:NH1	1:B:578:ASP:OD2	2.37	0.57
1:B:787:LYS:CE	3:B:1001:AE3:H1C2	2.25	0.57
1:A:256:LYS:HE3	2:A:1007:PGE:C3	2.33	0.57
1:A:497:CYS:HB3	1:A:820:HIS:CE1	2.39	0.57
1:B:604:ILE:HD13	1:B:631:THR:HG21	1.87	0.57
1:A:926:LYS:NZ	2:A:1010:PGE:O4	2.37	0.57
1:A:904:ARG:NH1	1:A:943:LYS:CB	2.68	0.57
1:B:188:LYS:HD3	1:B:189:ILE:H	1.71	0.56
1:B:499:GLN:NE2	1:B:817:SER:OG	2.35	0.56
1:B:915:TYR:HB2	1:B:921:LEU:HD12	1.88	0.56
1:A:322:SER:C	1:A:324:TRP:H	2.14	0.56
1:A:617:GLY:N	1:A:808:GLU:OE2	2.37	0.56
1:A:604:ILE:HG12	1:A:671:LYS:HD2	1.87	0.56
1:B:151:THR:HB	1:B:330:ALA:HB2	1.87	0.56
1:A:98:GLN:OE1	1:A:868:LYS:NZ	2.35	0.56
1:B:64:GLU:OE1	1:B:121:ASP:HB2	2.05	0.56
1:B:73:VAL:HG12	1:B:88:TRP:HB3	1.89	0.55
1:A:56:LYS:NZ	1:A:866:THR:O	2.40	0.55
1:A:496:ASP:OD2	2:A:1008:PGE:O4	2.24	0.55
1:A:145:TRP:O	1:A:147:ALA:N	2.33	0.54
1:B:352:LYS:HG3	1:B:353:PRO:HD2	1.89	0.54
1:B:245:ILE:HD11	1:B:287:TRP:CH2	2.43	0.54
1:A:139:LEU:HD12	1:A:145:TRP:CE2	2.43	0.54
1:A:760:GLY:O	1:A:764:THR:OG1	2.24	0.54
1:A:922:GLU:HB2	1:A:929:VAL:HB	1.88	0.53
1:A:256:LYS:NZ	2:A:1007:PGE:H42	2.23	0.53
1:A:268:GLN:NE2	1:A:280:TRP:CE3	2.69	0.53
1:A:814:GLY:O	1:A:816:ASN:N	2.41	0.53
1:B:73:VAL:HG13	1:B:87:VAL:HB	1.91	0.53
1:A:239:ARG:NH1	4:A:1107:HOH:O	2.41	0.53
1:B:118:ARG:NH2	1:B:126:GLU:OE2	2.42	0.52
1:B:270:LEU:HB2	1:B:280:TRP:CZ3	2.43	0.52
1:B:390:ARG:HE	1:B:442:LEU:HD13	1.73	0.52
1:A:674:LYS:O	1:A:676:LYS:N	2.42	0.52
1:A:491:MET:C	1:A:493:ILE:H	2.17	0.52
1:A:605:VAL:HG23	1:A:630:SER:HB3	1.92	0.52
1:A:103:GLU:OE2	2:A:1002:PGE:H42	2.10	0.52
1:B:86:ASP:OD1	1:B:86:ASP:N	2.43	0.52
1:B:145:TRP:O	1:B:268:GLN:NE2	2.43	0.52
1:A:206:ILE:HB	1:A:211:ILE:HD11	1.92	0.51
1:B:395:GLU:CD	1:B:499:GLN:HE21	2.18	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:491:MET:C	1:B:493:ILE:H	2.18	0.51
1:B:541:ASN:C	1:B:541:ASN:HD22	2.14	0.51
1:A:759:THR:HB	1:A:763:ALA:HB3	1.93	0.51
1:A:612:ASP:OD2	1:A:722:THR:OG1	2.24	0.51
1:A:180:ARG:HD3	1:A:321:ASP:OD1	2.11	0.51
1:A:404:ASN:OD1	1:A:811:THR:OG1	2.28	0.51
1:A:880:ASN:OD1	1:A:880:ASN:N	2.42	0.51
1:B:247:VAL:HG11	1:B:265:LEU:HD13	1.92	0.51
1:B:374:GLY:HA2	1:B:458:VAL:HB	1.93	0.51
1:A:566:ILE:HD11	1:A:637:ASP:OD2	2.10	0.51
1:A:280:TRP:CD1	1:A:280:TRP:N	2.77	0.51
1:B:633:TYR:HA	1:B:636:HIS:HB3	1.93	0.51
1:B:671:LYS:NZ	4:B:1105:HOH:O	2.44	0.50
1:B:722:THR:HG23	1:B:763:ALA:HB2	1.93	0.50
1:A:74:SER:HA	1:A:86:ASP:HB3	1.93	0.50
1:B:153:LYS:HG2	1:B:328:ILE:HG23	1.94	0.50
1:B:486:GLN:NE2	1:B:516:GLU:OE1	2.44	0.50
1:B:832:ARG:HB3	1:B:838:ARG:HG2	1.92	0.50
1:B:715:LEU:O	1:B:717:MET:N	2.45	0.50
1:B:38:GLU:OE2	1:B:48:LYS:HD3	2.12	0.50
1:B:169:LEU:HD22	1:B:257:GLU:HB2	1.94	0.50
1:B:352:LYS:HE2	1:B:451:GLU:HA	1.94	0.50
1:B:578:ASP:OD2	1:B:581:VAL:HB	2.12	0.50
1:B:797:GLU:OE2	3:B:1001:AE3:H2C1	2.12	0.49
1:A:49:LYS:HD3	1:A:140:MET:HE1	1.93	0.49
1:B:198:GLY:HA3	1:B:202:TYR:HB3	1.93	0.49
1:B:313:GLY:HA2	1:B:316:LEU:HD12	1.94	0.49
1:B:497:CYS:HB2	1:B:499:GLN:OE1	2.11	0.49
1:A:604:ILE:HD11	1:A:667:ALA:HB1	1.95	0.49
1:B:45:GLY:HA3	1:B:231:VAL:HG11	1.93	0.49
1:B:556:PRO:O	1:B:609:LYS:NZ	2.43	0.49
1:B:544:LEU:HD12	1:B:562:SER:HB3	1.95	0.48
1:B:874:ASN:HB2	1:B:885:ARG:HB3	1.95	0.48
1:B:105:LYS:NZ	4:B:1107:HOH:O	2.45	0.48
1:B:861:LYS:NZ	1:B:863:ARG:HH21	2.11	0.48
1:A:720:SER:H	1:A:723:LEU:HD12	1.78	0.47
1:B:477:LYS:HG2	2:B:1004:PGE:H22	1.96	0.47
1:A:112:VAL:HG22	1:A:136:GLU:HB2	1.96	0.47
1:A:149:TRP:HB3	1:A:330:ALA:HB3	1.96	0.47
1:B:202:TYR:CE1	1:B:217:ASP:HB3	2.49	0.47
1:B:759:THR:HB	1:B:763:ALA:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:509:ASP:N	1:A:509:ASP:OD1	2.48	0.47
1:B:74:SER:HA	1:B:86:ASP:OD1	2.14	0.47
1:B:508:GLY:H	1:B:561:TRP:CD1	2.32	0.47
1:B:545:SER:HB2	1:B:555:TYR:CZ	2.50	0.47
1:A:32:PRO:HG2	1:A:117:VAL:HG13	1.96	0.47
1:B:772:THR:HG23	1:B:777:ASP:HB3	1.97	0.47
1:B:511:GLN:HB2	1:B:564:ALA:HB2	1.96	0.46
1:A:359:PRO:HB2	1:A:402:THR:HG22	1.97	0.46
1:A:424:TYR:OH	1:A:426:GLU:OE1	2.29	0.46
1:B:385:SER:O	1:B:419:THR:OG1	2.26	0.46
1:B:188:LYS:CD	1:B:189:ILE:H	2.28	0.46
1:B:477:LYS:HD3	2:B:1004:PGE:H32	1.98	0.46
1:B:520:ASN:ND2	1:B:827:ASP:OD2	2.38	0.46
1:A:194:VAL:HG22	1:A:269:VAL:HG22	1.97	0.46
1:B:194:VAL:O	1:B:231:VAL:HA	2.16	0.46
1:A:68:ALA:HB3	1:A:120:TRP:HB2	1.98	0.46
1:A:351:ILE:HD11	1:A:372:TYR:CZ	2.51	0.46
1:B:236:LYS:HA	1:B:236:LYS:HD3	1.68	0.46
1:A:262:PHE:CZ	2:A:1007:PGE:H2	2.50	0.46
1:B:608:TYR:CD2	1:B:626:GLY:HA3	2.51	0.46
1:B:881:ILE:HD13	1:B:942:GLU:HA	1.97	0.46
1:A:253:ARG:HD2	1:A:551:TYR:CE2	2.51	0.45
1:B:213:ASP:OD1	1:B:213:ASP:N	2.47	0.45
1:B:312:LYS:HD3	1:B:313:GLY:H	1.80	0.45
1:B:474:LEU:HB2	3:B:1001:AE3:H5C1	1.98	0.45
1:B:602:ASP:OD2	1:B:602:ASP:N	2.45	0.45
1:B:872:GLU:HB2	1:B:887:SER:HB2	1.98	0.45
1:A:109:PRO:HA	1:A:140:MET:HG3	1.99	0.45
1:B:509:ASP:OD1	1:B:509:ASP:N	2.50	0.45
1:A:874:ASN:HB3	1:B:447:VAL:HG23	1.99	0.45
1:A:634:PHE:CD2	1:A:664:ILE:HD13	2.52	0.45
1:B:645:ALA:HB1	1:B:650:LYS:HB2	1.98	0.45
1:A:236:LYS:HA	1:A:236:LYS:HD2	1.79	0.45
1:A:382:THR:C	1:A:384:GLU:H	2.24	0.45
1:B:178:MET:HG2	1:B:248:ILE:HD13	1.98	0.45
1:B:27:SER:O	1:B:27:SER:OG	2.32	0.45
2:A:1004:PGE:H4	2:A:1004:PGE:H22	1.76	0.44
1:B:343:PRO:HA	1:B:344:PRO:HD3	1.92	0.44
1:A:140:MET:HB3	1:A:140:MET:HE3	1.72	0.44
1:B:194:VAL:HG23	1:B:234:VAL:HG21	2.00	0.44
1:B:312:LYS:HE2	1:B:312:LYS:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:717:MET:CG	1:A:752:ARG:HH22	2.20	0.44
1:A:357:TRP:CD2	2:A:1001:PGE:H3	2.53	0.44
1:A:756:HIS:HB3	1:A:790:PRO:HD2	1.99	0.44
1:B:217:ASP:CG	1:B:218:PRO:HA	2.42	0.44
1:A:74:SER:HB2	1:A:80:ILE:HB	1.99	0.44
1:B:625:SER:O	1:B:629:THR:HG23	2.18	0.44
1:B:59:HIS:CG	1:B:64:GLU:HG3	2.52	0.44
1:A:190:LYS:HD3	1:A:274:GLU:HA	1.99	0.44
1:A:523:MET:HE3	1:A:523:MET:HB3	1.79	0.44
1:A:592:TYR:HE2	1:A:637:ASP:OD1	2.01	0.44
1:A:923:GLU:HB3	1:A:928:ILE:HD13	1.99	0.43
1:B:924:ASN:C	1:B:926:LYS:H	2.25	0.43
1:A:152:LYS:HD3	1:A:177:PRO:HB3	2.00	0.43
1:A:924:ASN:ND2	4:A:1112:HOH:O	2.50	0.43
1:A:142:GLU:HA	1:A:145:TRP:HD1	1.83	0.43
1:A:908:LYS:HB2	1:A:908:LYS:HE3	1.65	0.43
1:A:908:LYS:HB2	1:A:940:TYR:HB2	2.01	0.43
1:A:926:LYS:CD	2:A:1010:PGE:H42	2.44	0.43
1:A:226:ARG:HH22	1:A:532:ASN:C	2.27	0.43
1:B:184:SER:C	1:B:185:ILE:CD1	2.92	0.43
1:A:87:VAL:HB	1:A:105:LYS:HB3	2.00	0.43
1:A:170:ASN:HD21	2:A:1007:PGE:H6	1.84	0.43
1:A:262:PHE:CZ	2:A:1007:PGE:C2	3.02	0.43
1:A:508:GLY:H	1:A:561:TRP:CD1	2.37	0.43
1:B:715:LEU:HD13	1:B:716:GLY:N	2.33	0.43
1:B:909:GLU:HB2	1:B:914:ILE:HD13	2.00	0.43
1:A:139:LEU:O	1:A:143:GLU:HB2	2.19	0.43
1:A:288:LYS:HG3	1:A:321:ASP:O	2.18	0.43
1:B:145:TRP:CE2	1:B:337:LEU:HD13	2.54	0.43
1:B:459:HIS:HA	1:B:491:MET:SD	2.59	0.43
1:A:604:ILE:N	1:A:604:ILE:CD1	2.82	0.43
1:A:635:TYR:HD2	1:A:730:LEU:HD13	1.83	0.43
1:B:177:PRO:HD2	1:B:249:LEU:HB2	2.01	0.43
1:B:613:TRP:CD1	1:B:614:CYS:HG	2.37	0.43
1:A:322:SER:C	1:A:324:TRP:N	2.76	0.42
1:B:56:LYS:HD3	1:B:869:GLY:HA2	2.02	0.42
1:B:604:ILE:HG21	1:B:671:LYS:HD2	2.00	0.42
1:B:822:MET:H	1:B:822:MET:HG3	1.59	0.42
1:A:377:LYS:NZ	1:A:454:GLU:OE1	2.52	0.42
1:A:805:GLU:OE2	2:A:1008:PGE:H62	2.20	0.42
1:B:288:LYS:HG3	1:B:321:ASP:OD2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:LYS:HD3	1:A:387:LYS:HA	1.97	0.42
1:A:545:SER:HB2	1:A:555:TYR:CZ	2.54	0.42
1:A:792:PHE:CE2	1:A:826:VAL:HG21	2.55	0.42
1:A:785:ASN:O	1:A:785:ASN:ND2	2.43	0.42
1:B:203:GLU:HB3	1:B:248:ILE:CG2	2.50	0.42
1:A:301:HIS:HA	1:A:413:ALA:H	1.84	0.42
1:B:301:HIS:HA	1:B:413:ALA:H	1.85	0.42
1:B:366:TYR:OH	1:B:448:PRO:HB2	2.19	0.42
1:A:42:ASN:HA	1:A:134:THR:O	2.19	0.42
1:B:479:HIS:ND1	1:B:827:ASP:OD1	2.32	0.42
1:A:561:TRP:CH2	1:A:761:ILE:HD11	2.55	0.42
1:A:668:PHE:HE1	1:A:727:PRO:HB3	1.84	0.41
1:B:59:HIS:CE1	1:B:64:GLU:OE2	2.73	0.41
1:B:491:MET:HB3	1:B:491:MET:HE3	1.76	0.41
1:B:861:LYS:HG3	1:B:874:ASN:OD1	2.19	0.41
1:B:924:ASN:O	1:B:925:GLU:HB3	2.20	0.41
1:A:898:VAL:HG12	1:A:941:MET:HE3	2.02	0.41
1:B:27:SER:C	1:B:29:ILE:CD1	2.91	0.41
1:B:674:LYS:O	1:B:676:LYS:N	2.51	0.41
1:B:898:VAL:HB	1:B:928:ILE:HB	2.02	0.41
1:A:881:ILE:HG12	1:A:942:GLU:HG2	2.03	0.41
1:B:207:ASN:ND2	1:B:237:ASN:O	2.50	0.41
1:B:373:THR:HG21	1:B:488:SER:HA	2.02	0.41
1:B:385:SER:O	1:B:387:LYS:N	2.46	0.41
1:A:447:VAL:HB	1:B:861:LYS:HB2	2.03	0.41
1:B:54:SER:HA	1:B:98:GLN:O	2.21	0.41
1:B:372:TYR:CD2	1:B:457:VAL:HG12	2.56	0.41
1:A:905:PHE:HA	1:A:942:GLU:O	2.19	0.41
1:A:109:PRO:HG3	1:A:140:MET:HB2	2.02	0.41
1:A:459:HIS:HA	1:A:491:MET:SD	2.60	0.41
1:A:874:ASN:O	1:A:884:MET:HA	2.21	0.41
1:B:35:LEU:HB3	1:B:115:TRP:CZ2	2.56	0.41
2:B:1005:PGE:H12	2:B:1005:PGE:H32	1.89	0.41
1:A:913:ILE:H	1:A:913:ILE:HD12	1.85	0.41
1:A:155:HIS:CE1	1:A:262:PHE:HB3	2.56	0.41
1:A:262:PHE:CE2	2:A:1007:PGE:C2	3.03	0.41
1:B:380:VAL:HG21	1:B:389:ILE:HD13	2.02	0.41
1:B:509:ASP:OD2	1:B:822:MET:HG2	2.20	0.41
1:B:677:ALA:HA	1:B:741:VAL:HG22	2.02	0.41
1:B:750:ILE:HG12	1:B:755:TYR:CZ	2.56	0.41
1:B:366:TYR:CE1	1:B:453:ILE:HG13	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:792:PHE:CE2	1:B:826:VAL:HG21	2.56	0.40
1:A:239:ARG:HG2	1:A:240:ASP:H	1.86	0.40
3:A:1005:AE3:H5C1	3:A:1005:AE3:H3C1	1.68	0.40
1:B:145:TRP:HA	1:B:268:GLN:HE22	1.85	0.40
1:B:349:LYS:HB2	1:B:349:LYS:HE2	1.97	0.40
1:A:116:LYS:HG2	1:A:132:VAL:HG22	2.04	0.40
1:A:290:THR:OG1	1:A:291:TYR:N	2.55	0.40
1:B:180:ARG:HD3	1:B:321:ASP:HB3	2.02	0.40
1:B:839:VAL:HG12	4:B:1134:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	879/944 (93%)	818 (93%)	54 (6%)	7 (1%)	16	29
1	B	882/944 (93%)	814 (92%)	63 (7%)	5 (1%)	21	36
All	All	1761/1888 (93%)	1632 (93%)	117 (7%)	12 (1%)	18	32

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	323	THR
1	A	241	GLY
1	A	322	SER
1	A	815	MET
1	B	716	GLY
1	A	498	PRO
1	B	140	MET
1	B	673	LEU
1	A	320	ASP

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Mol	Chain	Res	Type
1	A	907	VAL
1	B	386	GLY
1	B	166	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	777/840 (92%)	767 (99%)	10 (1%)	61	75
1	B	778/840 (93%)	772 (99%)	6 (1%)	73	83
All	All	1555/1680 (93%)	1539 (99%)	16 (1%)	68	80

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	139	LEU
1	A	256	LYS
1	A	280	TRP
1	A	655	LYS
1	A	662	ASP
1	A	717	MET
1	A	904	ARG
1	A	908	LYS
1	A	917	LYS
1	A	941	MET
1	B	26	SER
1	B	191	ARG
1	B	680	SER
1	B	779	VAL
1	B	787	LYS
1	B	809	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	ASN
1	A	133	ASN
1	A	237	ASN
1	A	404	ASN
1	A	731	ASN
1	B	63	ASN
1	B	268	GLN
1	B	486	GLN
1	B	541	ASN
1	B	786	GLN
1	B	816	ASN
1	B	876	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

16 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	AE3	B	1002	-	7,7,8	0.32	0	6,6,7	0.24	0
2	PGE	B	1004	-	9,9,9	0.32	0	8,8,8	0.29	0
2	PGE	A	1001	-	9,9,9	0.33	0	8,8,8	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PGE	A	1003	-	9,9,9	0.32	0	8,8,8	0.29	0
2	PGE	A	1004	-	9,9,9	0.31	0	8,8,8	0.30	0
3	AE3	A	1005	-	8,8,8	0.38	0	7,7,7	0.55	0
3	AE3	A	1006	-	7,7,8	0.31	0	6,6,7	0.27	0
3	AE3	B	1001	-	8,8,8	0.37	0	7,7,7	0.28	0
2	PGE	B	1003	-	9,9,9	0.32	0	8,8,8	0.25	0
2	PGE	B	1005	-	9,9,9	0.32	0	8,8,8	0.29	0
3	AE3	B	1006	-	5,5,8	0.37	0	4,4,7	0.29	0
2	PGE	A	1007	-	9,9,9	0.28	0	8,8,8	0.63	0
2	PGE	A	1009	-	9,9,9	0.32	0	8,8,8	0.32	0
2	PGE	A	1010	-	9,9,9	0.51	0	8,8,8	0.79	0
2	PGE	A	1008	-	9,9,9	0.32	0	8,8,8	0.26	0
2	PGE	A	1002	-	9,9,9	0.32	0	8,8,8	0.28	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
3	AE3	B	1002	-	-	2/5/5/6	-
2	PGE	B	1004	-	-	6/7/7/7	-
2	PGE	A	1001	-	-	1/7/7/7	-
2	PGE	A	1003	-	-	0/7/7/7	-
2	PGE	A	1004	-	-	3/7/7/7	-
3	AE3	A	1005	-	-	4/6/6/6	-
3	AE3	A	1006	-	-	0/5/5/6	-
3	AE3	B	1001	-	-	2/6/6/6	-
2	PGE	B	1003	-	-	1/7/7/7	-
2	PGE	B	1005	-	-	5/7/7/7	-
3	AE3	B	1006	-	-	1/3/3/6	-
2	PGE	A	1007	-	-	6/7/7/7	-
2	PGE	A	1009	-	-	0/7/7/7	-
2	PGE	A	1010	-	-	4/7/7/7	-
2	PGE	A	1008	-	-	2/7/7/7	-
2	PGE	A	1002	-	-	2/7/7/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1010	PGE	C4-C3-O2-C2
3	A	1005	AE3	C3-C4-O3-C5
3	A	1005	AE3	O2-C3-C4-O3
2	B	1005	PGE	O2-C3-C4-O3
2	A	1004	PGE	O2-C3-C4-O3
3	B	1002	AE3	O2-C3-C4-O3
2	A	1007	PGE	O2-C3-C4-O3
2	A	1010	PGE	O3-C5-C6-O4
2	B	1004	PGE	O1-C1-C2-O2
2	B	1004	PGE	O2-C3-C4-O3
2	A	1010	PGE	O1-C1-C2-O2
2	B	1003	PGE	O3-C5-C6-O4
2	A	1008	PGE	O2-C3-C4-O3
2	A	1010	PGE	O2-C3-C4-O3
2	A	1008	PGE	C6-C5-O3-C4
2	A	1007	PGE	C1-C2-O2-C3
3	B	1001	AE3	C6-C5-O3-C4
2	A	1004	PGE	C6-C5-O3-C4
2	A	1001	PGE	C3-C4-O3-C5
2	B	1004	PGE	O3-C5-C6-O4
2	A	1002	PGE	C3-C4-O3-C5
2	B	1005	PGE	O1-C1-C2-O2
2	B	1004	PGE	C3-C4-O3-C5
3	A	1005	AE3	C6-C5-O3-C4
2	A	1007	PGE	C4-C3-O2-C2
3	A	1005	AE3	C1-C2-O2-C3
2	B	1004	PGE	C1-C2-O2-C3
2	A	1007	PGE	O3-C5-C6-O4
3	B	1001	AE3	C3-C4-O3-C5
2	A	1004	PGE	C4-C3-O2-C2
2	B	1004	PGE	C4-C3-O2-C2
2	B	1005	PGE	C4-C3-O2-C2
2	A	1007	PGE	C6-C5-O3-C4
2	A	1007	PGE	C3-C4-O3-C5
2	A	1002	PGE	O3-C5-C6-O4
2	B	1005	PGE	C3-C4-O3-C5
2	B	1005	PGE	C1-C2-O2-C3
3	B	1006	AE3	C1-C2-O2-C3
3	B	1002	AE3	C3-C4-O3-C5

There are no ring outliers.

10 monomers are involved in 41 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1004	PGE	5	0
2	A	1001	PGE	8	0
2	A	1004	PGE	1	0
3	A	1005	AE3	1	0
3	B	1001	AE3	4	0
2	B	1005	PGE	1	0
2	A	1007	PGE	12	0
2	A	1010	PGE	6	0
2	A	1008	PGE	2	0
2	A	1002	PGE	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	883/944 (93%)	0.58	30 (3%) 48 46	42, 58, 81, 102	0
1	B	886/944 (93%)	0.56	33 (3%) 45 43	42, 60, 83, 97	0
All	All	1769/1888 (93%)	0.57	63 (3%) 46 44	42, 59, 82, 102	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	814	GLY	4.2
1	B	714	SER	4.0
1	B	715	LEU	4.0
1	A	84	ILE	3.7
1	B	212	GLY	3.7
1	B	213	ASP	3.6
1	B	541	ASN	3.5
1	A	719	THR	3.4
1	B	911	ASP	3.3
1	B	84	ILE	3.3
1	A	717	MET	3.2
1	B	716	GLY	3.2
1	A	141	ASN	3.2
1	B	638	VAL	3.1
1	A	85	GLY	3.0
1	A	165	ALA	3.0
1	B	25	LYS	3.0
1	A	904	ARG	2.9
1	A	935	GLY	2.9
1	B	276	ASP	2.9
1	A	943	LYS	2.9
1	B	274	GLU	2.9
1	A	154	GLU	2.8
1	A	718	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	188	LYS	2.7
1	A	166	PRO	2.7
1	B	939	PHE	2.6
1	B	598	LYS	2.5
1	A	86	ASP	2.5
1	B	717	MET	2.5
1	A	323	THR	2.4
1	A	680	SER	2.4
1	A	25	LYS	2.4
1	A	26	SER	2.4
1	A	911	ASP	2.3
1	A	144	ASN	2.3
1	A	245	ILE	2.2
1	A	604	ILE	2.2
1	B	788	THR	2.2
1	A	731	ASN	2.2
1	B	736	ASP	2.2
1	A	316	LEU	2.2
1	A	238	ILE	2.2
1	A	123	LYS	2.2
1	B	26	SER	2.2
1	A	281	ILE	2.2
1	B	672	PHE	2.2
1	B	235	SER	2.2
1	A	888	VAL	2.1
1	B	238	ILE	2.1
1	A	925	GLU	2.1
1	B	83	ASN	2.1
1	B	415	ASP	2.1
1	B	627	GLU	2.1
1	B	669	ASN	2.1
1	B	313	GLY	2.1
1	B	799	GLY	2.1
1	A	188	LYS	2.0
1	B	278	ILE	2.0
1	B	886	ILE	2.0
1	B	607	PHE	2.0
1	B	169	LEU	2.0
1	B	843	GLY	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	AE3	B	1006	6/9	0.56	0.33	65,68,82,86	0
3	AE3	B	1001	9/9	0.60	0.28	60,78,100,105	0
2	PGE	A	1007	10/10	0.73	0.19	79,82,87,88	0
3	AE3	A	1005	9/9	0.73	0.32	76,108,124,125	0
2	PGE	A	1004	10/10	0.76	0.18	64,77,83,84	0
3	AE3	B	1002	8/9	0.79	0.17	59,65,67,69	0
2	PGE	B	1004	10/10	0.79	0.18	64,71,78,80	0
3	AE3	A	1006	8/9	0.83	0.14	67,75,77,80	0
2	PGE	B	1005	10/10	0.85	0.16	55,62,67,67	0
2	PGE	A	1010	10/10	0.85	0.20	20,20,20,20	0
2	PGE	A	1008	10/10	0.85	0.16	55,62,67,69	0
2	PGE	A	1003	10/10	0.86	0.15	53,62,85,92	0
2	PGE	A	1001	10/10	0.88	0.13	61,67,72,73	0
2	PGE	A	1002	10/10	0.88	0.14	62,66,74,75	0
2	PGE	B	1003	10/10	0.89	0.12	66,73,86,87	0
2	PGE	A	1009	10/10	0.90	0.13	57,63,67,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.