



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:30 PM GMT

PDB ID : 1TZ7
Title : Aquifex aeolicus amylomaltase
Authors : Barends, T.R.M.; Korf, H.; Kaper, T.; van der Maarel, M.J.E.C.; Dijkhuizen, L.; Dijkstra, B.W.
Deposited on : 2004-07-09
Resolution : 2.15 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

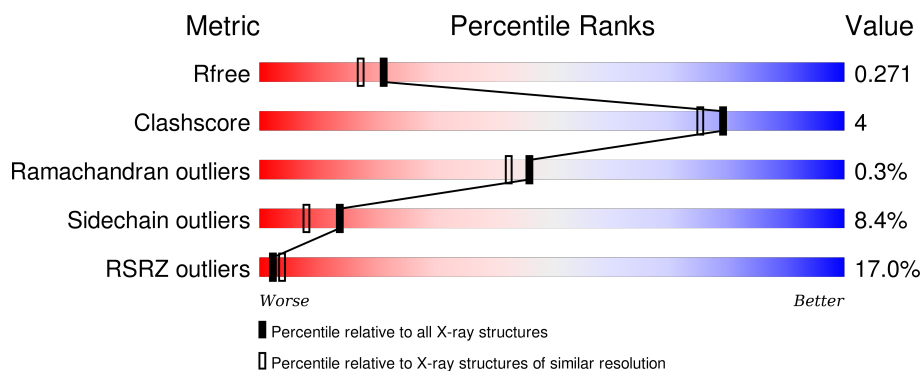
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	1045 (2.16-2.16)
Clashscore	102246	1152 (2.16-2.16)
Ramachandran outliers	100387	1131 (2.16-2.16)
Sidechain outliers	100360	1131 (2.16-2.16)
RSRZ outliers	91569	1050 (2.16-2.16)

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	505	<div> <div>14%</div> <div>81%</div> <div>14%</div> <div>• •</div> </div>
1	B	505	<div> <div>19%</div> <div>80%</div> <div>14%</div> <div>• •</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
2	MPD	A	3282	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8508 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 4-alpha-glucanotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	0	0
			4147	2714	695	732	6			
1	B	487	Total	C	N	O	S	0	0	0
			4143	2712	694	731	6			

There are 42 discrepancies between the modelled and reference sequences:

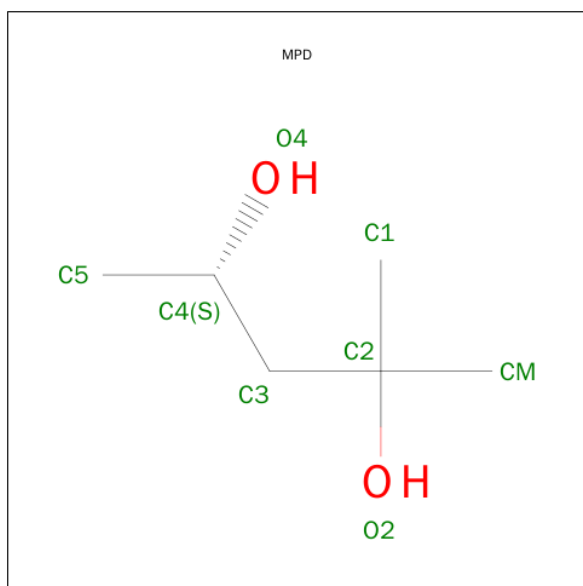
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP O66937
A	-18	GLY	-	EXPRESSION TAG	UNP O66937
A	-17	SER	-	EXPRESSION TAG	UNP O66937
A	-16	SER	-	EXPRESSION TAG	UNP O66937
A	-15	HIS	-	EXPRESSION TAG	UNP O66937
A	-14	HIS	-	EXPRESSION TAG	UNP O66937
A	-13	HIS	-	EXPRESSION TAG	UNP O66937
A	-12	HIS	-	EXPRESSION TAG	UNP O66937
A	-11	HIS	-	EXPRESSION TAG	UNP O66937
A	-10	HIS	-	EXPRESSION TAG	UNP O66937
A	-9	SER	-	EXPRESSION TAG	UNP O66937
A	-8	SER	-	EXPRESSION TAG	UNP O66937
A	-7	GLY	-	EXPRESSION TAG	UNP O66937
A	-6	LEU	-	EXPRESSION TAG	UNP O66937
A	-5	VAL	-	EXPRESSION TAG	UNP O66937
A	-4	PRO	-	EXPRESSION TAG	UNP O66937
A	-3	ARG	-	EXPRESSION TAG	UNP O66937
A	-2	GLY	-	EXPRESSION TAG	UNP O66937
A	-1	SER	-	EXPRESSION TAG	UNP O66937
A	0	HIS	-	EXPRESSION TAG	UNP O66937
A	274	LEU	HIS	CONFLICT	UNP O66937
B	-19	MET	-	EXPRESSION TAG	UNP O66937
B	-18	GLY	-	EXPRESSION TAG	UNP O66937
B	-17	SER	-	EXPRESSION TAG	UNP O66937
B	-16	SER	-	EXPRESSION TAG	UNP O66937

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	HIS	-	EXPRESSION TAG	UNP O66937
B	-14	HIS	-	EXPRESSION TAG	UNP O66937
B	-13	HIS	-	EXPRESSION TAG	UNP O66937
B	-12	HIS	-	EXPRESSION TAG	UNP O66937
B	-11	HIS	-	EXPRESSION TAG	UNP O66937
B	-10	HIS	-	EXPRESSION TAG	UNP O66937
B	-9	SER	-	EXPRESSION TAG	UNP O66937
B	-8	SER	-	EXPRESSION TAG	UNP O66937
B	-7	GLY	-	EXPRESSION TAG	UNP O66937
B	-6	LEU	-	EXPRESSION TAG	UNP O66937
B	-5	VAL	-	EXPRESSION TAG	UNP O66937
B	-4	PRO	-	EXPRESSION TAG	UNP O66937
B	-3	ARG	-	EXPRESSION TAG	UNP O66937
B	-2	GLY	-	EXPRESSION TAG	UNP O66937
B	-1	SER	-	EXPRESSION TAG	UNP O66937
B	0	HIS	-	EXPRESSION TAG	UNP O66937
B	274	LEU	HIS	CONFLICT	UNP O66937

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	6	2		

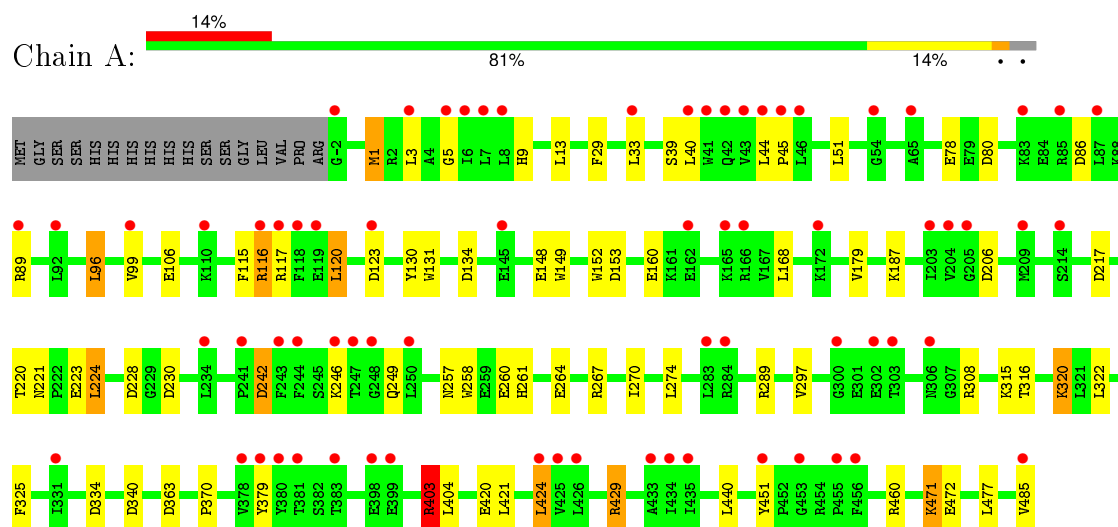
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	98	Total 98	O 98	0	0
3	B	112	Total 112	O 112	0	0

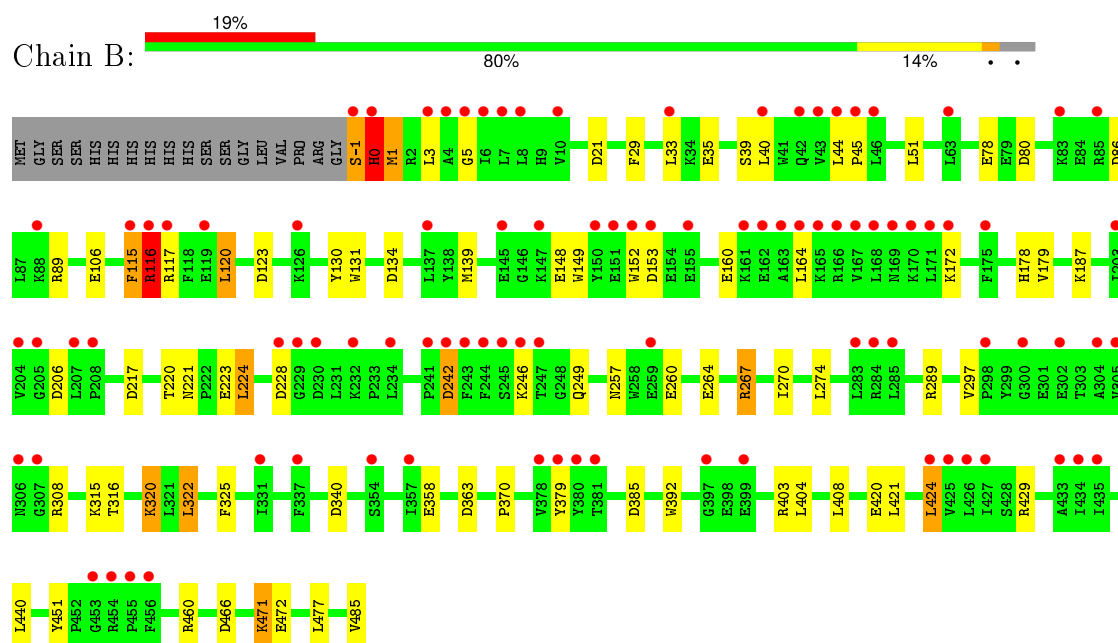
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: 4-alpha-glucanotransferase



• Molecule 1: 4-alpha-glucanotransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	156.75Å 156.75Å 169.16Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.15 29.62 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.8 (15.00-2.15) 99.8 (29.62-2.15)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.99 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.215 , 0.243 0.260 , 0.271	Depositor DCC
R_{free} test set	3374 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	37.7	Xtriage
Anisotropy	0.378	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	1 of 66778 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8508	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.94% 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

5.1 Standard geometry

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

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	0.45	0/4273	0.68	24/5768 (0.4%)
1	B	0.46	0/4269	0.68	25/5763 (0.4%)
All	All	0.45	0/8542	0.68	49/11531 (0.4%)

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	1
All	All	0	2

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	403	ARG	NE-CZ-NH1	12.77	126.69	120.30
1	A	403	ARG	NE-CZ-NH2	-12.12	114.24	120.30
1	B	403	ARG	NE-CZ-NH1	-12.09	114.26	120.30
1	B	403	ARG	NE-CZ-NH2	10.92	125.76	120.30
1	B	429	ARG	NE-CZ-NH1	-10.06	115.27	120.30
1	B	429	ARG	NE-CZ-NH2	10.04	125.32	120.30
1	B	460	ARG	NE-CZ-NH2	-9.27	115.67	120.30
1	A	267	ARG	NE-CZ-NH2	-9.24	115.68	120.30
1	A	460	ARG	NE-CZ-NH2	9.17	124.88	120.30
1	A	308	ARG	NE-CZ-NH2	-9.13	115.73	120.30
1	B	267	ARG	NE-CZ-NH1	-9.12	115.74	120.30
1	B	267	ARG	NE-CZ-NH2	9.10	124.85	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	460	ARG	NE-CZ-NH1	-9.09	115.75	120.30
1	B	460	ARG	NE-CZ-NH1	9.08	124.84	120.30
1	A	267	ARG	NE-CZ-NH1	9.05	124.82	120.30
1	B	308	ARG	NE-CZ-NH1	-8.98	115.81	120.30
1	A	429	ARG	NE-CZ-NH1	8.97	124.78	120.30
1	A	429	ARG	NE-CZ-NH2	-8.87	115.86	120.30
1	A	308	ARG	NE-CZ-NH1	8.84	124.72	120.30
1	B	308	ARG	NE-CZ-NH2	8.76	124.68	120.30
1	A	403	ARG	CD-NE-CZ	6.36	132.51	123.60
1	A	403	ARG	CG-CD-NE	5.64	123.64	111.80
1	B	228	ASP	CB-CG-OD2	5.55	123.30	118.30
1	A	228	ASP	CB-CG-OD2	5.46	123.21	118.30
1	A	206	ASP	CB-CG-OD2	5.41	123.17	118.30
1	B	340	ASP	CB-CG-OD2	5.39	123.15	118.30
1	B	217	ASP	CB-CG-OD2	5.37	123.14	118.30
1	B	206	ASP	CB-CG-OD2	5.36	123.12	118.30
1	B	134	ASP	CB-CG-OD2	5.31	123.08	118.30
1	B	172	LYS	CB-CA-C	5.31	121.01	110.40
1	A	340	ASP	CB-CG-OD2	5.29	123.06	118.30
1	A	334	ASP	CB-CG-OD2	5.29	123.06	118.30
1	B	363	ASP	CB-CG-OD2	5.28	123.05	118.30
1	A	363	ASP	CB-CG-OD2	5.27	123.04	118.30
1	B	80	ASP	CB-CG-OD2	5.23	123.00	118.30
1	A	153	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	80	ASP	CB-CG-OD2	5.21	122.98	118.30
1	A	230	ASP	CB-CG-OD2	5.18	122.96	118.30
1	B	86	ASP	CB-CG-OD2	5.17	122.96	118.30
1	B	385	ASP	CB-CG-OD2	5.17	122.96	118.30
1	A	86	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	217	ASP	CB-CG-OD2	5.12	122.91	118.30
1	B	123	ASP	CB-CG-OD2	5.11	122.90	118.30
1	A	123	ASP	CB-CG-OD2	5.11	122.90	118.30
1	B	153	ASP	CB-CG-OD2	5.10	122.89	118.30
1	B	466	ASP	CB-CG-OD2	5.10	122.89	118.30
1	B	21	ASP	CB-CG-OD2	5.08	122.87	118.30
1	B	403	ARG	CG-CD-NE	-5.01	101.27	111.80
1	A	134	ASP	CB-CG-OD2	5.01	122.81	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	115	PHE	Peptide
1	B	115	PHE	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4147	0	4071	27	0
1	B	4143	0	4068	31	0
2	A	8	0	14	0	0
3	A	98	0	0	1	0
3	B	112	0	0	1	0
All	All	8508	0	8153	58	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 4.

All (58) 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:B:-1:SER:HA	1:B:0:HIS:CB	1.90	1.01
1:B:-1:SER:HA	1:B:0:HIS:HB2	1.41	0.99
1:A:224:LEU:HG	1:A:261:HIS:CE1	2.04	0.92
1:B:-1:SER:HA	1:B:0:HIS:CG	2.07	0.88
1:B:-1:SER:CA	1:B:0:HIS:HB2	2.10	0.82
1:B:257:ASN:HD22	1:B:260:GLU:HB2	1.50	0.77
1:A:257:ASN:HD22	1:A:260:GLU:HB2	1.51	0.75
1:B:29:PHE:CZ	1:B:33:LEU:HD21	2.28	0.68
1:A:29:PHE:CZ	1:A:33:LEU:HD21	2.31	0.65
1:A:3:LEU:HD23	1:A:39:SER:CB	2.26	0.65
1:B:3:LEU:HD23	1:B:39:SER:CB	2.27	0.64
1:A:29:PHE:O	1:A:33:LEU:HD23	1.99	0.63
1:A:3:LEU:HD23	1:A:39:SER:HB2	1.83	0.61
1:A:242:ASP:N	1:A:242:ASP:OD1	2.35	0.59
1:B:242:ASP:N	1:B:242:ASP:OD1	2.35	0.58
1:B:3:LEU:HD23	1:B:39:SER:HB2	1.85	0.58
1:B:29:PHE:O	1:B:33:LEU:HD23	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:TRP:HA	1:A:261:HIS:HD2	1.68	0.58
1:A:270:ILE:HD11	1:A:320:LYS:HD3	1.91	0.53
1:B:270:ILE:HD11	1:B:320:LYS:HD3	1.90	0.52
1:A:96:LEU:HB3	1:A:99:VAL:HG13	1.92	0.51
1:A:370:PRO:HG2	1:A:424:LEU:HG	1.92	0.51
1:A:485:VAL:HG11	3:A:3288:HOH:O	2.10	0.51
1:A:221:ASN:HB3	1:A:224:LEU:HD22	1.93	0.50
1:B:370:PRO:HG2	1:B:424:LEU:HG	1.93	0.49
1:B:139:MET:HB3	1:B:178:HIS:CE1	2.47	0.49
1:B:149:TRP:HA	1:B:152:TRP:CD2	2.47	0.49
1:A:5:GLY:HA3	1:A:40:LEU:HB2	1.94	0.49
1:B:5:GLY:HA3	1:B:40:LEU:HB2	1.94	0.49
1:B:115:PHE:HA	1:B:116:ARG:HB2	1.95	0.48
1:B:221:ASN:HB3	1:B:224:LEU:HD22	1.94	0.47
1:A:274:LEU:HD12	1:A:325:PHE:HE2	1.80	0.47
1:A:149:TRP:HA	1:A:152:TRP:CD2	2.49	0.47
1:B:130:TYR:CE1	1:B:131:TRP:HD1	2.33	0.46
1:A:471:LYS:HG3	1:A:472:GLU:N	2.30	0.46
1:B:485:VAL:HG11	3:B:498:HOH:O	2.15	0.46
1:A:130:TYR:CE1	1:A:131:TRP:HD1	2.34	0.46
1:A:1:MET:O	1:A:1:MET:HG3	2.16	0.46
1:A:404:LEU:HD11	1:A:421:LEU:HD21	1.97	0.45
1:B:44:LEU:HB3	1:B:45:PRO:HD2	1.98	0.45
1:A:44:LEU:HB3	1:A:45:PRO:HD2	1.99	0.45
1:B:404:LEU:HD11	1:B:421:LEU:HD21	1.97	0.45
1:A:224:LEU:HG	1:A:261:HIS:ND1	2.32	0.45
1:B:44:LEU:HB3	1:B:45:PRO:CD	2.47	0.44
1:A:9:HIS:CE1	1:A:44:LEU:HB2	2.52	0.44
1:B:471:LYS:HG3	1:B:472:GLU:N	2.32	0.44
1:B:1:MET:HG3	1:B:1:MET:O	2.16	0.44
1:A:44:LEU:HB3	1:A:45:PRO:CD	2.48	0.43
1:B:29:PHE:CE2	1:B:33:LEU:HD21	2.53	0.43
1:A:403:ARG:HH11	1:A:403:ARG:HG3	1.83	0.43
1:B:274:LEU:HD12	1:B:325:PHE:HE2	1.84	0.43
1:B:117:ARG:HG2	1:B:120:LEU:HD23	2.01	0.42
1:B:322:LEU:HA	1:B:322:LEU:HD13	1.94	0.42
1:A:258:TRP:HA	1:A:261:HIS:CD2	2.50	0.41
1:A:117:ARG:HG2	1:A:120:LEU:HD12	2.01	0.41
1:B:-1:SER:CB	1:B:0:HIS:HB2	2.51	0.40
1:B:257:ASN:ND2	1:B:260:GLU:HB2	2.25	0.40
1:B:358:GLU:HG3	1:B:392:TRP:CD2	2.56	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	486/505 (96%)	475 (98%)	10 (2%)	1 (0%)	52	51
1	B	485/505 (96%)	474 (98%)	9 (2%)	2 (0%)	39	34
All	All	971/1010 (96%)	949 (98%)	19 (2%)	3 (0%)	46	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	0	HIS
1	B	116	ARG
1	A	116	ARG

5.3.2 Protein sidechains [i](#)

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	436/451 (97%)	400 (92%)	36 (8%)	14	8
1	B	436/451 (97%)	399 (92%)	37 (8%)	13	8
All	All	872/902 (97%)	799 (92%)	73 (8%)	14	8

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	13	LEU
1	A	51	LEU
1	A	78	GLU
1	A	89	ARG
1	A	96	LEU
1	A	106	GLU
1	A	116	ARG
1	A	120	LEU
1	A	148	GLU
1	A	160	GLU
1	A	168	LEU
1	A	179	VAL
1	A	187	LYS
1	A	220	THR
1	A	223	GLU
1	A	224	LEU
1	A	242	ASP
1	A	246	LYS
1	A	249	GLN
1	A	264	GLU
1	A	289	ARG
1	A	297	VAL
1	A	315	LYS
1	A	316	THR
1	A	320	LYS
1	A	322	LEU
1	A	379	TYR
1	A	403	ARG
1	A	420	GLU
1	A	424	LEU
1	A	429	ARG
1	A	440	LEU
1	A	451	TYR
1	A	471	LYS
1	A	477	LEU
1	B	-1	SER
1	B	0	HIS
1	B	1	MET
1	B	35	GLU
1	B	51	LEU
1	B	78	GLU
1	B	89	ARG

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Mol	Chain	Res	Type
1	B	106	GLU
1	B	116	ARG
1	B	120	LEU
1	B	148	GLU
1	B	160	GLU
1	B	164	LEU
1	B	179	VAL
1	B	187	LYS
1	B	220	THR
1	B	223	GLU
1	B	224	LEU
1	B	242	ASP
1	B	246	LYS
1	B	249	GLN
1	B	264	GLU
1	B	267	ARG
1	B	289	ARG
1	B	297	VAL
1	B	315	LYS
1	B	316	THR
1	B	320	LYS
1	B	322	LEU
1	B	379	TYR
1	B	408	LEU
1	B	420	GLU
1	B	424	LEU
1	B	440	LEU
1	B	451	TYR
1	B	471	LYS
1	B	477	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	42	GLN
1	A	55	ASN
1	A	249	GLN
1	A	257	ASN
1	A	436	GLN
1	B	9	HIS
1	B	42	GLN

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Mol	Chain	Res	Type
1	B	55	ASN
1	B	249	GLN
1	B	257	ASN
1	B	384	HIS
1	B	436	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	MPD	A	3282	-	6,7,7	0.33	0	7,10,10	0.41	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MPD	A	3282	-	-	0/5/5/5	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [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	488/505 (96%)	0.87	71 (14%) 3 5	21, 32, 56, 107	0
1	B	487/505 (96%)	1.02	95 (19%) 1 2	16, 32, 56, 98	0
All	All	975/1010 (96%)	0.94	166 (17%) 2 4	16, 32, 56, 107	0

All (166) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	243	PHE	15.1
1	B	244	PHE	10.6
1	B	455	PRO	8.4
1	A	243	PHE	8.4
1	B	246	LYS	7.4
1	A	244	PHE	7.2
1	A	246	LYS	6.9
1	B	247	THR	6.0
1	B	116	ARG	5.7
1	A	116	ARG	5.7
1	B	166	ARG	5.4
1	B	456	PHE	5.1
1	B	300	GLY	5.0
1	A	426	LEU	5.0
1	B	162	GLU	4.9
1	B	453	GLY	4.8
1	B	150	TYR	4.7
1	B	-1	SER	4.7
1	B	306	ASN	4.6
1	A	85	ARG	4.6
1	B	434	ILE	4.5
1	A	302	GLU	4.5
1	A	380	TYR	4.5
1	B	163	ALA	4.4

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Mol	Chain	Res	Type	RSRZ
1	B	43	VAL	4.4
1	A	8	LEU	4.4
1	A	456	PHE	4.3
1	B	172	LYS	4.3
1	B	5	GLY	4.2
1	A	6	ILE	4.1
1	B	7	LEU	4.0
1	A	453	GLY	4.0
1	A	162	GLU	3.9
1	B	153	ASP	3.9
1	B	380	TYR	3.9
1	A	434	ILE	3.9
1	A	46	LEU	3.9
1	B	46	LEU	3.9
1	B	229	GLY	3.8
1	A	41	TRP	3.8
1	B	242	ASP	3.7
1	B	171	LEU	3.7
1	A	425	VAL	3.6
1	B	305	VAL	3.6
1	B	204	VAL	3.6
1	A	247	THR	3.5
1	A	435	ILE	3.5
1	B	298	PRO	3.5
1	A	300	GLY	3.5
1	A	166	ARG	3.5
1	A	44	LEU	3.4
1	A	5	GLY	3.4
1	B	302	GLU	3.4
1	A	43	VAL	3.4
1	B	145	GLU	3.4
1	A	204	VAL	3.3
1	A	7	LEU	3.3
1	A	241	PRO	3.3
1	B	234	LEU	3.3
1	B	115	PHE	3.2
1	B	378	VAL	3.2
1	B	207	LEU	3.2
1	B	435	ILE	3.2
1	B	44	LEU	3.2
1	A	306	ASN	3.2
1	B	433	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	165	LYS	3.2
1	A	379	TYR	3.2
1	B	3	LEU	3.2
1	A	455	PRO	3.2
1	A	433	ALA	3.1
1	B	45	PRO	3.1
1	A	-2	GLY	3.1
1	B	85	ARG	3.0
1	A	378	VAL	3.0
1	A	45	PRO	2.9
1	B	283	LEU	2.9
1	A	145	GLU	2.9
1	B	425	VAL	2.9
1	B	167	VAL	2.9
1	B	126	LYS	2.8
1	B	381	THR	2.8
1	B	245	SER	2.8
1	B	304	ALA	2.8
1	A	119	GLU	2.8
1	A	381	THR	2.8
1	B	232	LYS	2.7
1	B	168	LEU	2.7
1	B	426	LEU	2.7
1	A	248	GLY	2.7
1	B	307	GLY	2.7
1	B	6	ILE	2.7
1	B	379	TYR	2.6
1	B	284	ARG	2.6
1	B	83	LYS	2.6
1	A	205	GLY	2.6
1	B	337	PHE	2.6
1	A	33	LEU	2.6
1	B	40	LEU	2.6
1	B	137	LEU	2.6
1	B	208	PRO	2.6
1	A	203	ILE	2.5
1	B	241	PRO	2.5
1	B	454	ARG	2.5
1	A	3	LEU	2.5
1	B	119	GLU	2.5
1	A	214	SER	2.5
1	A	118	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	92	LEU	2.5
1	A	165	LYS	2.5
1	A	110	LYS	2.5
1	A	40	LEU	2.5
1	B	427	ILE	2.5
1	B	147	LYS	2.5
1	B	8	LEU	2.4
1	B	164	LEU	2.4
1	B	424	LEU	2.4
1	A	399	GLU	2.4
1	B	155	GLU	2.4
1	B	169	ASN	2.4
1	B	203	ILE	2.4
1	B	230	ASP	2.4
1	B	205	GLY	2.4
1	A	424	LEU	2.3
1	B	4	ALA	2.3
1	A	117	ARG	2.3
1	B	117	ARG	2.3
1	A	83	LYS	2.3
1	B	354	SER	2.3
1	A	89	ARG	2.3
1	B	152	TRP	2.3
1	A	284	ARG	2.3
1	A	283	LEU	2.3
1	B	0	HIS	2.2
1	A	250	LEU	2.2
1	A	451	TYR	2.2
1	A	383	THR	2.2
1	B	175	PHE	2.2
1	B	399	GLU	2.2
1	A	54	GLY	2.2
1	A	172	LYS	2.2
1	A	234	LEU	2.1
1	A	303	THR	2.1
1	A	65	ALA	2.1
1	B	10	VAL	2.1
1	A	398	GLU	2.1
1	B	151	GLU	2.1
1	B	331	ILE	2.1
1	A	123	ASP	2.1
1	B	285	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	42	GLN	2.1
1	B	259	GLU	2.1
1	A	331	ILE	2.1
1	A	87	LEU	2.1
1	B	63	LEU	2.1
1	A	99	VAL	2.1
1	A	485	VAL	2.1
1	B	228	ASP	2.1
1	B	161	LYS	2.0
1	B	170	LYS	2.0
1	B	397	GLY	2.0
1	A	209	MET	2.0
1	B	357	ILE	2.0
1	B	33	LEU	2.0
1	A	42	GLN	2.0
1	B	88	LYS	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
2	MPD	A	3282	8/8	0.91	0.36	4.23	57,65,66,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.