



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:38 AM GMT

PDB ID : 2B4K
Title : Acetobacter turbidans alpha-amino acid ester hydrolase complexed with phenylglycine
Authors : Barends, T.R.M.; Polderman-Tijmes, J.J.; Jekel, P.A.; Williams, C.; Wybenga, G.; Janssen, D.B.; Dijkstra, B.W.
Deposited on : 2005-09-26
Resolution : 3.30 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

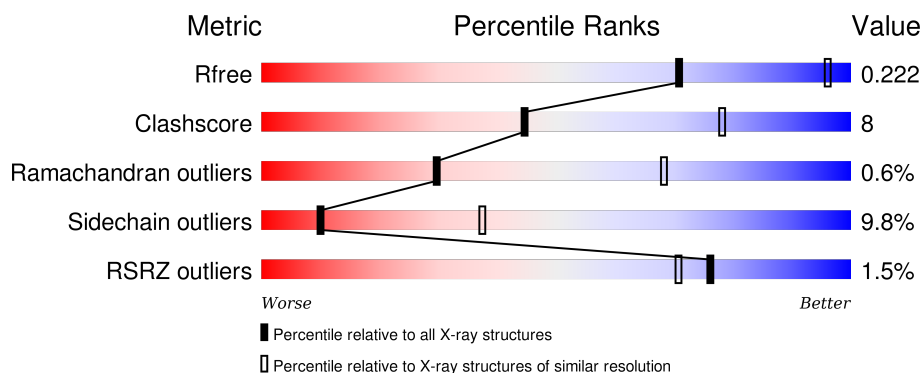
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	2060 (3.40-3.20)
Clashscore	102246	1058 (3.38-3.22)
Ramachandran outliers	100387	1038 (3.38-3.22)
Sidechain outliers	100360	1037 (3.38-3.22)
RSRZ outliers	91569	2070 (3.40-3.20)

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	652	<div> <div>2%</div> <div>69% 23% 5%</div> </div>
1	B	652	<div> <div>2%</div> <div>68% 24% 5%</div> </div>
1	C	652	<div> <div>2%</div> <div>66% 25% 5%</div> </div>
1	D	652	<div> <div>2%</div> <div>68% 24% 5%</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	PG9	A	3600	-	-	-	X
2	PG9	B	3599	-	-	-	X
2	PG9	D	3598	-	-	-	X
3	GOL	B	2432	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 19594 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-amino acid ester hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	617	Total	C	N	O	S	0	0	0
			4883	3111	845	907	20			
1	B	617	Total	C	N	O	S	0	0	0
			4883	3111	845	907	20			
1	C	617	Total	C	N	O	S	0	0	0
			4883	3111	845	907	20			
1	D	617	Total	C	N	O	S	0	0	0
			4883	3111	845	907	20			

There are 100 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	668	LYS	-	EXPRESSION TAG	UNP Q8VRK8
A	669	LEU	-	EXPRESSION TAG	UNP Q8VRK8
A	670	GLY	-	EXPRESSION TAG	UNP Q8VRK8
A	671	PRO	-	EXPRESSION TAG	UNP Q8VRK8
A	672	GLU	-	myc tag	UNP Q8VRK8
A	673	GLN	-	myc tag	UNP Q8VRK8
A	674	LYS	-	myc tag	UNP Q8VRK8
A	675	LEU	-	myc tag	UNP Q8VRK8
A	676	ILE	-	myc tag	UNP Q8VRK8
A	677	SER	-	myc tag	UNP Q8VRK8
A	678	GLU	-	myc tag	UNP Q8VRK8
A	679	GLU	-	myc tag	UNP Q8VRK8
A	680	ASP	-	myc tag	UNP Q8VRK8
A	681	LEU	-	myc tag	UNP Q8VRK8
A	682	ASN	-	EXPRESSION TAG	UNP Q8VRK8
A	683	SER	-	EXPRESSION TAG	UNP Q8VRK8
A	684	ALA	-	EXPRESSION TAG	UNP Q8VRK8
A	685	VAL	-	EXPRESSION TAG	UNP Q8VRK8
A	686	ASP	-	EXPRESSION TAG	UNP Q8VRK8
A	687	HIS	-	EXPRESSION TAG	UNP Q8VRK8
A	688	HIS	-	EXPRESSION TAG	UNP Q8VRK8

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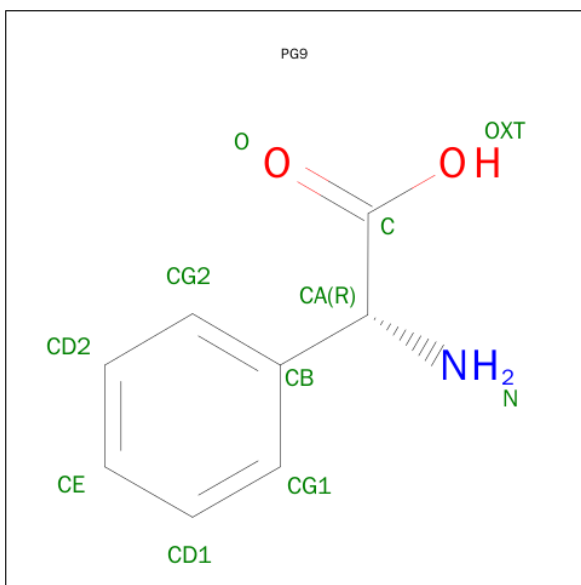
Chain	Residue	Modelled	Actual	Comment	Reference
A	689	HIS	-	EXPRESSION TAG	UNP Q8VRK8
A	690	HIS	-	EXPRESSION TAG	UNP Q8VRK8
A	691	HIS	-	EXPRESSION TAG	UNP Q8VRK8
A	692	HIS	-	EXPRESSION TAG	UNP Q8VRK8
B	668	LYS	-	EXPRESSION TAG	UNP Q8VRK8
B	669	LEU	-	EXPRESSION TAG	UNP Q8VRK8
B	670	GLY	-	EXPRESSION TAG	UNP Q8VRK8
B	671	PRO	-	EXPRESSION TAG	UNP Q8VRK8
B	672	GLU	-	myc tag	UNP Q8VRK8
B	673	GLN	-	myc tag	UNP Q8VRK8
B	674	LYS	-	myc tag	UNP Q8VRK8
B	675	LEU	-	myc tag	UNP Q8VRK8
B	676	ILE	-	myc tag	UNP Q8VRK8
B	677	SER	-	myc tag	UNP Q8VRK8
B	678	GLU	-	myc tag	UNP Q8VRK8
B	679	GLU	-	myc tag	UNP Q8VRK8
B	680	ASP	-	myc tag	UNP Q8VRK8
B	681	LEU	-	myc tag	UNP Q8VRK8
B	682	ASN	-	EXPRESSION TAG	UNP Q8VRK8
B	683	SER	-	EXPRESSION TAG	UNP Q8VRK8
B	684	ALA	-	EXPRESSION TAG	UNP Q8VRK8
B	685	VAL	-	EXPRESSION TAG	UNP Q8VRK8
B	686	ASP	-	EXPRESSION TAG	UNP Q8VRK8
B	687	HIS	-	EXPRESSION TAG	UNP Q8VRK8
B	688	HIS	-	EXPRESSION TAG	UNP Q8VRK8
B	689	HIS	-	EXPRESSION TAG	UNP Q8VRK8
B	690	HIS	-	EXPRESSION TAG	UNP Q8VRK8
B	691	HIS	-	EXPRESSION TAG	UNP Q8VRK8
B	692	HIS	-	EXPRESSION TAG	UNP Q8VRK8
C	668	LYS	-	EXPRESSION TAG	UNP Q8VRK8
C	669	LEU	-	EXPRESSION TAG	UNP Q8VRK8
C	670	GLY	-	EXPRESSION TAG	UNP Q8VRK8
C	671	PRO	-	EXPRESSION TAG	UNP Q8VRK8
C	672	GLU	-	myc tag	UNP Q8VRK8
C	673	GLN	-	myc tag	UNP Q8VRK8
C	674	LYS	-	myc tag	UNP Q8VRK8
C	675	LEU	-	myc tag	UNP Q8VRK8
C	676	ILE	-	myc tag	UNP Q8VRK8
C	677	SER	-	myc tag	UNP Q8VRK8
C	678	GLU	-	myc tag	UNP Q8VRK8
C	679	GLU	-	myc tag	UNP Q8VRK8
C	680	ASP	-	myc tag	UNP Q8VRK8

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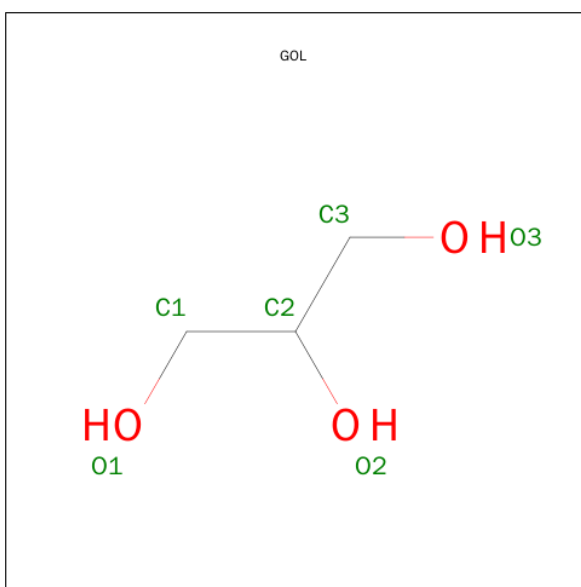
Chain	Residue	Modelled	Actual	Comment	Reference
C	681	LEU	-	myc tag	UNP Q8VRK8
C	682	ASN	-	EXPRESSION TAG	UNP Q8VRK8
C	683	SER	-	EXPRESSION TAG	UNP Q8VRK8
C	684	ALA	-	EXPRESSION TAG	UNP Q8VRK8
C	685	VAL	-	EXPRESSION TAG	UNP Q8VRK8
C	686	ASP	-	EXPRESSION TAG	UNP Q8VRK8
C	687	HIS	-	EXPRESSION TAG	UNP Q8VRK8
C	688	HIS	-	EXPRESSION TAG	UNP Q8VRK8
C	689	HIS	-	EXPRESSION TAG	UNP Q8VRK8
C	690	HIS	-	EXPRESSION TAG	UNP Q8VRK8
C	691	HIS	-	EXPRESSION TAG	UNP Q8VRK8
C	692	HIS	-	EXPRESSION TAG	UNP Q8VRK8
D	668	LYS	-	EXPRESSION TAG	UNP Q8VRK8
D	669	LEU	-	EXPRESSION TAG	UNP Q8VRK8
D	670	GLY	-	EXPRESSION TAG	UNP Q8VRK8
D	671	PRO	-	EXPRESSION TAG	UNP Q8VRK8
D	672	GLU	-	myc tag	UNP Q8VRK8
D	673	GLN	-	myc tag	UNP Q8VRK8
D	674	LYS	-	myc tag	UNP Q8VRK8
D	675	LEU	-	myc tag	UNP Q8VRK8
D	676	ILE	-	myc tag	UNP Q8VRK8
D	677	SER	-	myc tag	UNP Q8VRK8
D	678	GLU	-	myc tag	UNP Q8VRK8
D	679	GLU	-	myc tag	UNP Q8VRK8
D	680	ASP	-	myc tag	UNP Q8VRK8
D	681	LEU	-	myc tag	UNP Q8VRK8
D	682	ASN	-	EXPRESSION TAG	UNP Q8VRK8
D	683	SER	-	EXPRESSION TAG	UNP Q8VRK8
D	684	ALA	-	EXPRESSION TAG	UNP Q8VRK8
D	685	VAL	-	EXPRESSION TAG	UNP Q8VRK8
D	686	ASP	-	EXPRESSION TAG	UNP Q8VRK8
D	687	HIS	-	EXPRESSION TAG	UNP Q8VRK8
D	688	HIS	-	EXPRESSION TAG	UNP Q8VRK8
D	689	HIS	-	EXPRESSION TAG	UNP Q8VRK8
D	690	HIS	-	EXPRESSION TAG	UNP Q8VRK8
D	691	HIS	-	EXPRESSION TAG	UNP Q8VRK8
D	692	HIS	-	EXPRESSION TAG	UNP Q8VRK8

- Molecule 2 is D-PHENYLGLYCINE (three-letter code: PG9) (formula: $C_8H_9NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			11	8	1	2		
2	B	1	Total	C	N	O	0	0
			11	8	1	2		
2	A	1	Total	C	N	O	0	0
			11	8	1	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

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

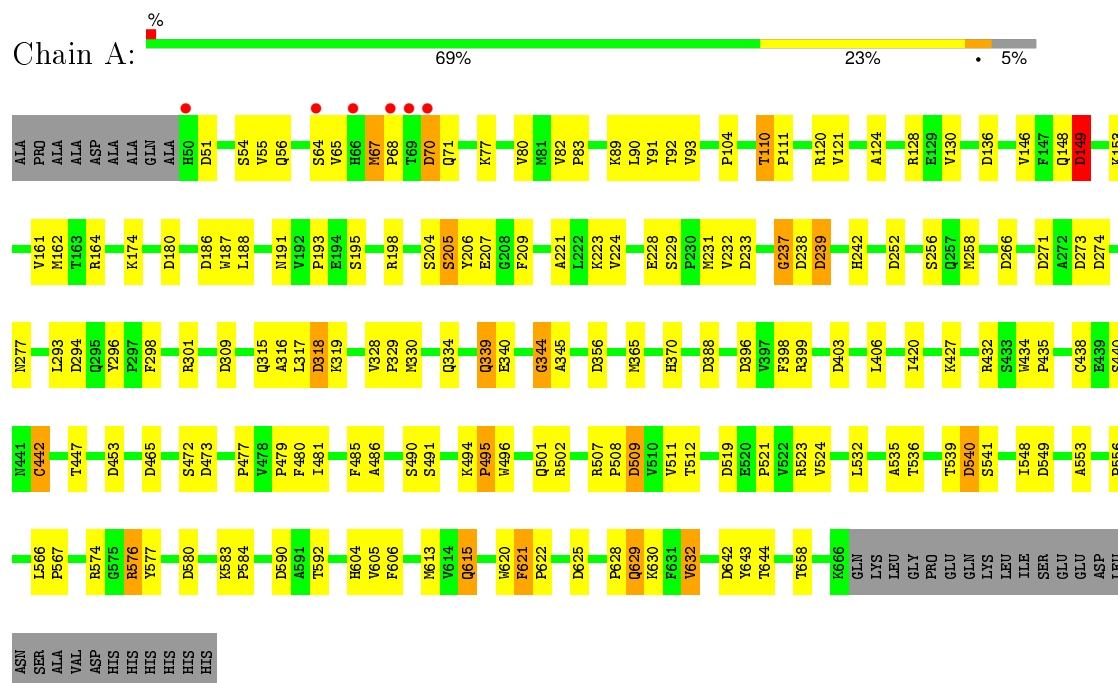
- Molecule 4 is water.

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

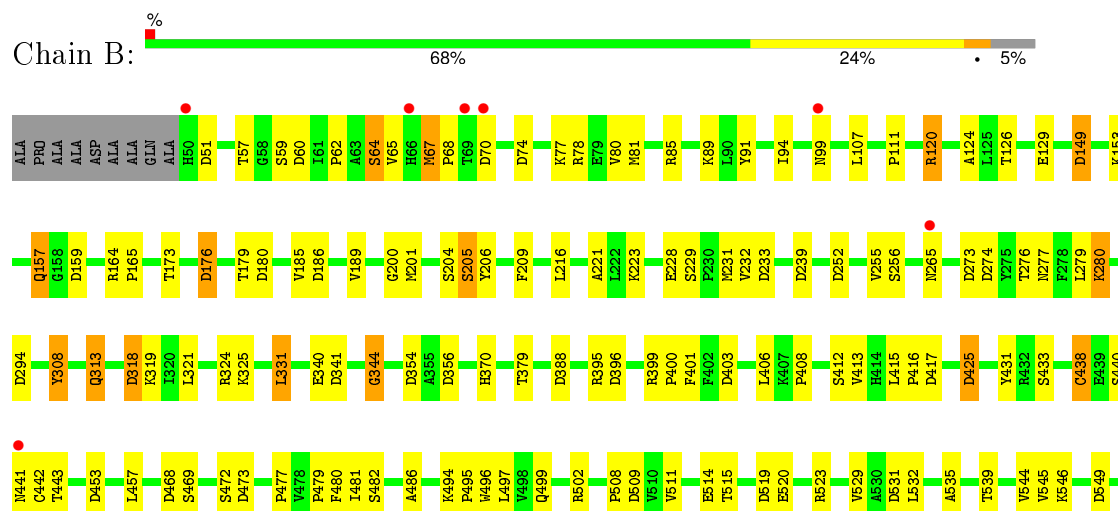
3 Residue-property plots

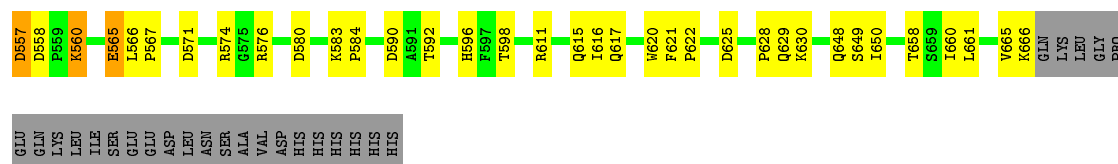
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: alpha-amino acid ester hydrolase

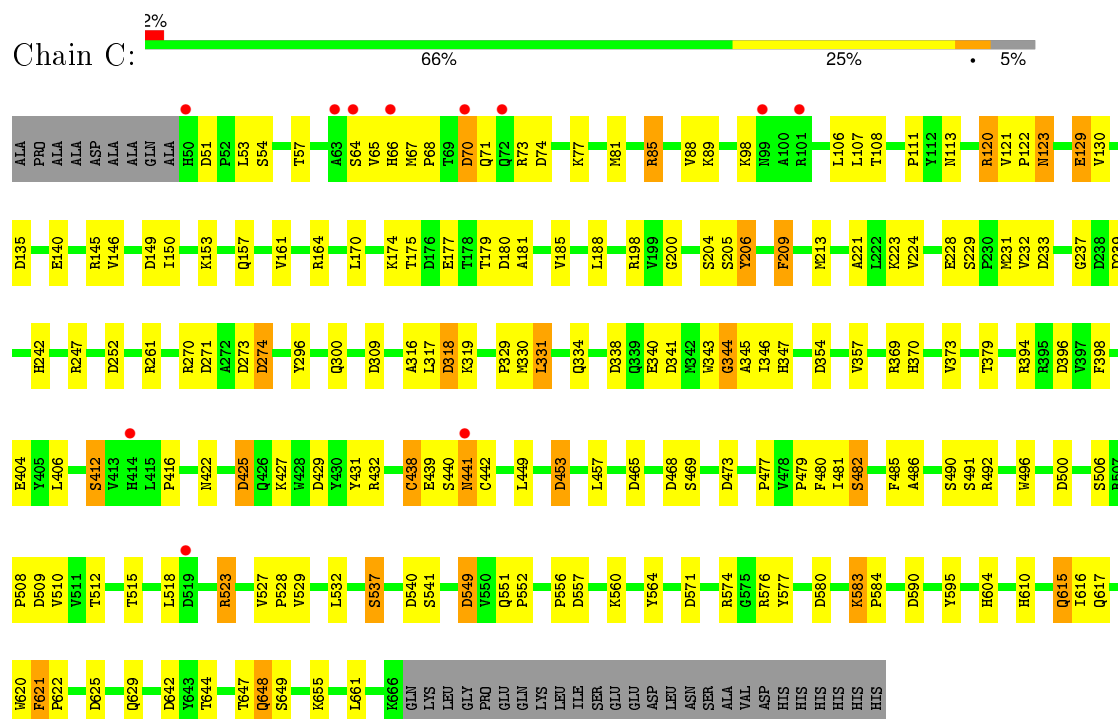


- Molecule 1: alpha-amino acid ester hydrolase

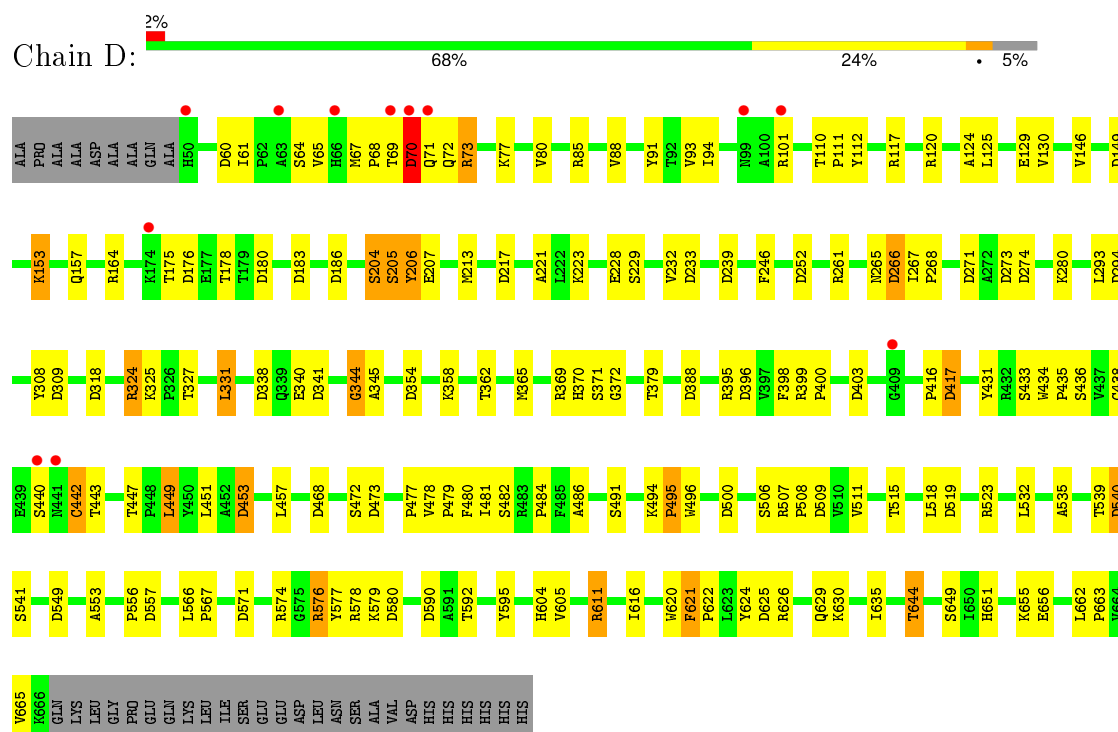




• Molecule 1: alpha-amino acid ester hydrolase



• Molecule 1: alpha-amino acid ester hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	340.91Å 340.91Å 340.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 3.30 19.95 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.4 (15.00-3.30) 99.4 (19.95-3.30)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 3.29Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.258 , 0.289 0.220 , 0.222	Depositor DCC
R_{free} test set	4687 reflections (5.10%)	DCC
Wilson B-factor (Å ²)	48.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , -6.8	EDS
Estimated twinning fraction	0.000 for -l,-k,-h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	0 of 97300 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	19594	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.59% 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: GOL, PG9

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	0.41	0/5042	0.73	28/6887 (0.4%)
1	B	0.42	0/5042	0.73	32/6887 (0.5%)
1	C	0.43	0/5042	0.73	31/6887 (0.5%)
1	D	0.42	0/5042	0.73	33/6887 (0.5%)
All	All	0.42	0/20168	0.73	124/27548 (0.5%)

There are no bond length outliers.

All (124) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	580	ASP	CB-CG-OD2	6.14	123.83	118.30
1	D	233	ASP	CB-CG-OD2	6.09	123.78	118.30
1	B	509	ASP	CB-CG-OD2	6.04	123.74	118.30
1	C	473	ASP	CB-CG-OD2	6.00	123.70	118.30
1	C	233	ASP	CB-CG-OD2	5.97	123.67	118.30
1	C	74	ASP	CB-CG-OD2	5.95	123.65	118.30
1	B	74	ASP	CB-CG-OD2	5.95	123.65	118.30
1	C	51	ASP	CB-CG-OD2	5.91	123.62	118.30
1	D	580	ASP	CB-CG-OD2	5.89	123.60	118.30
1	D	473	ASP	CB-CG-OD2	5.88	123.59	118.30
1	C	509	ASP	CB-CG-OD2	5.87	123.58	118.30
1	B	356	ASP	CB-CG-OD2	5.84	123.56	118.30
1	C	580	ASP	CB-CG-OD2	5.84	123.55	118.30
1	A	473	ASP	CB-CG-OD2	5.81	123.53	118.30
1	A	580	ASP	CB-CG-OD2	5.80	123.52	118.30
1	B	625	ASP	CB-CG-OD2	5.80	123.52	118.30
1	D	509	ASP	CB-CG-OD2	5.79	123.51	118.30
1	B	233	ASP	CB-CG-OD2	5.75	123.47	118.30
1	A	318	ASP	CB-CG-OD2	5.73	123.45	118.30
1	C	540	ASP	CB-CG-OD2	5.72	123.44	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	549	ASP	CB-CG-OD2	5.71	123.44	118.30
1	B	549	ASP	CB-CG-OD2	5.71	123.44	118.30
1	C	642	ASP	CB-CG-OD2	5.69	123.42	118.30
1	A	625	ASP	CB-CG-OD2	5.68	123.41	118.30
1	A	51	ASP	CB-CG-OD2	5.66	123.39	118.30
1	B	403	ASP	CB-CG-OD2	5.64	123.38	118.30
1	D	625	ASP	CB-CG-OD2	5.64	123.37	118.30
1	A	273	ASP	CB-CG-OD2	5.63	123.37	118.30
1	C	549	ASP	CB-CG-OD2	5.62	123.36	118.30
1	B	473	ASP	CB-CG-OD2	5.62	123.36	118.30
1	D	252	ASP	CB-CG-OD2	5.60	123.34	118.30
1	A	266	ASP	CB-CG-OD2	5.59	123.33	118.30
1	C	625	ASP	CB-CG-OD2	5.57	123.31	118.30
1	B	557	ASP	CB-CG-OD2	5.52	123.27	118.30
1	B	149	ASP	CB-CG-OD2	5.52	123.27	118.30
1	C	318	ASP	CB-CG-OD2	5.52	123.27	118.30
1	A	180	ASP	CB-CG-OD2	5.50	123.25	118.30
1	C	252	ASP	CB-CG-OD2	5.48	123.23	118.30
1	C	453	ASP	CB-CG-OD2	5.47	123.22	118.30
1	D	266	ASP	CB-CG-OD2	5.46	123.21	118.30
1	D	453	ASP	CB-CG-OD2	5.42	123.18	118.30
1	A	642	ASP	CB-CG-OD2	5.42	123.17	118.30
1	C	70	ASP	CB-CG-OD2	5.42	123.17	118.30
1	D	217	ASP	CB-CG-OD2	5.42	123.17	118.30
1	D	354	ASP	CB-CG-OD2	5.41	123.17	118.30
1	C	571	ASP	CB-CG-OD2	5.41	123.17	118.30
1	B	51	ASP	CB-CG-OD2	5.40	123.16	118.30
1	D	403	ASP	CB-CG-OD2	5.39	123.15	118.30
1	B	531	ASP	CB-CG-OD2	5.38	123.15	118.30
1	D	70	ASP	CB-CG-OD2	5.38	123.14	118.30
1	C	273	ASP	CB-CG-OD2	5.36	123.13	118.30
1	A	239	ASP	CB-CG-OD2	5.36	123.12	118.30
1	B	252	ASP	CB-CG-OD2	5.36	123.12	118.30
1	B	453	ASP	CB-CG-OD2	5.35	123.11	118.30
1	D	341	ASP	CB-CG-OD2	5.34	123.11	118.30
1	A	309	ASP	CB-CG-OD2	5.34	123.10	118.30
1	A	509	ASP	CB-CG-OD2	5.34	123.10	118.30
1	D	500	ASP	CB-CG-OD2	5.33	123.10	118.30
1	D	540	ASP	CB-CG-OD2	5.33	123.09	118.30
1	A	238	ASP	CB-CG-OD2	5.32	123.09	118.30
1	B	388	ASP	CB-CG-OD2	5.32	123.09	118.30
1	D	273	ASP	CB-CG-OD2	5.30	123.07	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	180	ASP	CB-CG-OD2	5.30	123.07	118.30
1	B	590	ASP	CB-CG-OD2	5.30	123.07	118.30
1	A	549	ASP	CB-CG-OD2	5.29	123.06	118.30
1	D	318	ASP	CB-CG-OD2	5.29	123.06	118.30
1	A	540	ASP	CB-CG-OD2	5.29	123.06	118.30
1	C	341	ASP	CB-CG-OD2	5.29	123.06	118.30
1	A	396	ASP	CB-CG-OD2	5.29	123.06	118.30
1	D	590	ASP	CB-CG-OD2	5.28	123.05	118.30
1	C	135	ASP	CB-CG-OD2	5.27	123.05	118.30
1	C	239	ASP	CB-CG-OD2	5.27	123.05	118.30
1	D	180	ASP	CB-CG-OD2	5.23	123.01	118.30
1	A	233	ASP	CB-CG-OD2	5.23	123.00	118.30
1	A	388	ASP	CB-CG-OD2	5.22	123.00	118.30
1	D	468	ASP	CB-CG-OD2	5.21	122.99	118.30
1	B	180	ASP	CB-CG-OD2	5.20	122.97	118.30
1	D	396	ASP	CB-CG-OD2	5.20	122.98	118.30
1	B	186	ASP	CB-CG-OD2	5.19	122.97	118.30
1	C	354	ASP	CB-CG-OD2	5.19	122.97	118.30
1	C	309	ASP	CB-CG-OD2	5.19	122.97	118.30
1	B	341	ASP	CB-CG-OD2	5.19	122.97	118.30
1	A	356	ASP	CB-CG-OD2	5.18	122.96	118.30
1	D	388	ASP	CB-CG-OD2	5.18	122.96	118.30
1	A	590	ASP	CB-CG-OD2	5.17	122.96	118.30
1	C	274	ASP	CB-CG-OD2	5.17	122.96	118.30
1	D	239	ASP	CB-CG-OD2	5.17	122.96	118.30
1	D	149	ASP	CB-CG-OD2	5.17	122.95	118.30
1	B	273	ASP	CB-CG-OD2	5.16	122.94	118.30
1	D	557	ASP	CB-CG-OD2	5.16	122.94	118.30
1	A	149	ASP	CB-CG-OD2	5.16	122.94	118.30
1	D	571	ASP	CB-CG-OD2	5.16	122.94	118.30
1	A	519	ASP	CB-CG-OD2	5.15	122.94	118.30
1	D	186	ASP	CB-CG-OD2	5.15	122.93	118.30
1	B	468	ASP	CB-CG-OD2	5.15	122.93	118.30
1	C	590	ASP	CB-CG-OD2	5.14	122.93	118.30
1	B	354	ASP	CB-CG-OD2	5.14	122.92	118.30
1	B	294	ASP	CB-CG-OD2	5.13	122.92	118.30
1	D	60	ASP	CB-CG-OD2	5.13	122.91	118.30
1	D	309	ASP	CB-CG-OD2	5.12	122.91	118.30
1	B	60	ASP	CB-CG-OD2	5.12	122.91	118.30
1	C	149	ASP	CB-CG-OD2	5.11	122.90	118.30
1	B	176	ASP	CB-CG-OD2	5.09	122.88	118.30
1	B	425	ASP	CB-CG-OD2	5.09	122.88	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	396	ASP	CB-CG-OD2	5.08	122.87	118.30
1	A	186	ASP	CB-CG-OD2	5.08	122.87	118.30
1	A	252	ASP	CB-CG-OD2	5.07	122.87	118.30
1	A	465	ASP	CB-CG-OD2	5.07	122.87	118.30
1	A	70	ASP	CB-CG-OD2	5.07	122.86	118.30
1	C	425	ASP	CB-CG-OD2	5.07	122.86	118.30
1	C	468	ASP	CB-CG-OD2	5.06	122.86	118.30
1	C	557	ASP	CB-CG-OD2	5.06	122.85	118.30
1	D	176	ASP	CB-CG-OD2	5.06	122.85	118.30
1	A	294	ASP	CB-CG-OD2	5.05	122.85	118.30
1	B	571	ASP	CB-CG-OD2	5.05	122.85	118.30
1	B	239	ASP	CB-CG-OD2	5.05	122.84	118.30
1	C	271	ASP	CB-CG-OD2	5.05	122.84	118.30
1	A	403	ASP	CB-CG-OD2	5.04	122.84	118.30
1	C	500	ASP	CB-CG-OD2	5.03	122.83	118.30
1	B	318	ASP	CB-CG-OD2	5.03	122.83	118.30
1	D	294	ASP	CB-CG-OD2	5.03	122.83	118.30
1	D	338	ASP	CB-CG-OD2	5.02	122.82	118.30
1	C	396	ASP	CB-CG-OD2	5.01	122.81	118.30
1	B	159	ASP	CB-CG-OD2	5.01	122.81	118.30

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	4883	0	4648	74	0
1	B	4883	0	4648	77	0
1	C	4883	0	4648	77	0
1	D	4883	0	4648	80	0
2	A	11	0	8	2	0
2	B	11	0	8	1	0
2	D	11	0	8	4	0
3	A	12	0	16	1	0
3	B	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	6	0	8	0	0
4	A	1	0	0	0	0
4	B	4	0	0	0	0
All	All	19594	0	18648	298	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 8.

All (298) 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:438:CYS:HG	1:A:442:CYS:HG	1.03	0.94
1:D:232:VAL:HB	1:D:344:GLY:HA2	1.56	0.87
1:C:70:ASP:HA	1:C:73:ARG:HD2	1.57	0.85
1:D:205:SER:OG	1:D:206:TYR:N	2.09	0.84
1:C:482:SER:O	1:C:492:ARG:NH1	2.11	0.83
1:B:232:VAL:HB	1:B:344:GLY:HA2	1.66	0.78
1:B:205:SER:OG	1:B:370:HIS:NE2	2.19	0.76
1:A:501:GLN:HG2	1:A:567:PRO:HG3	1.69	0.72
1:A:232:VAL:HB	1:A:344:GLY:HA2	1.75	0.69
1:D:532:LEU:HB3	1:D:595:TYR:HB2	1.75	0.67
1:B:274:ASP:HA	1:B:277:ASN:HB2	1.77	0.66
1:C:532:LEU:HB3	1:C:595:TYR:HB2	1.77	0.66
1:D:438:CYS:SG	1:D:442:CYS:HB2	2.36	0.66
1:D:438:CYS:HG	1:D:442:CYS:HG	1.43	0.65
1:B:457:LEU:HB2	1:B:515:THR:HG22	1.79	0.64
1:C:440:SER:O	1:C:441:ASN:HB2	1.97	0.64
1:C:481:ILE:HG21	1:C:496:TRP:HB2	1.80	0.63
1:C:161:VAL:HB	1:C:164:ARG:HG3	1.81	0.62
1:D:205:SER:HG	1:D:206:TYR:H	1.46	0.62
1:D:449:LEU:HD11	1:D:457:LEU:HB3	1.83	0.61
1:B:438:CYS:CB	1:B:442:CYS:HG	2.13	0.61
1:D:578:ARG:HH21	1:D:579:LYS:HE2	1.66	0.60
1:A:80:VAL:HG11	1:A:187:TRP:CZ2	2.37	0.60
1:B:438:CYS:HB3	1:B:442:CYS:HG	1.67	0.59
1:A:494:LYS:HB2	1:A:495:PRO:HD3	1.83	0.59
1:D:651:HIS:HD2	1:D:656:GLU:O	1.85	0.59
1:B:438:CYS:HG	1:B:442:CYS:HG	1.48	0.59
1:A:621:PHE:CG	1:A:622:PRO:HA	2.37	0.59
1:B:438:CYS:SG	1:B:442:CYS:SG	3.01	0.58
1:C:438:CYS:HB3	1:C:442:CYS:HG	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:438:CYS:HG	1:A:442:CYS:CB	2.16	0.58
1:B:340:GLU:HB2	1:B:370:HIS:CG	2.38	0.58
1:B:440:SER:O	1:B:441:ASN:HB2	2.03	0.57
1:D:399:ARG:HB3	1:D:400:PRO:HD3	1.86	0.57
1:A:477:PRO:HB2	1:A:621:PHE:HB2	1.85	0.57
1:B:438:CYS:HB3	1:B:442:CYS:SG	2.45	0.57
1:A:204:SER:HA	1:A:228:GLU:O	2.05	0.57
1:B:204:SER:HA	1:B:228:GLU:O	2.05	0.57
1:A:111:PRO:HG3	1:A:206:TYR:CG	2.40	0.57
1:D:204:SER:HA	1:D:228:GLU:O	2.05	0.56
1:C:479:PRO:HA	1:C:621:PHE:O	2.05	0.56
1:D:274:ASP:HB2	1:D:486:ALA:HA	1.88	0.56
1:A:556:PRO:HB2	1:B:120:ARG:CZ	2.36	0.56
1:C:438:CYS:SG	1:C:439:GLU:N	2.78	0.56
1:D:574:ARG:HB3	1:D:577:TYR:HD1	1.71	0.56
1:A:576:ARG:NH1	1:A:632:VAL:O	2.34	0.56
1:A:553:ALA:HB1	1:D:271:ASP:HA	1.88	0.55
1:D:232:VAL:HB	1:D:344:GLY:CA	2.33	0.55
1:B:176:ASP:OD1	1:B:179:THR:HG23	2.06	0.55
1:D:91:TYR:HE2	1:D:93:VAL:CG2	2.19	0.55
1:C:209:PHE:HB2	1:C:231:MET:HB2	1.88	0.55
1:C:512:THR:HG23	1:C:615:GLN:HG3	1.88	0.55
1:A:274:ASP:HA	1:A:277:ASN:HB2	1.88	0.55
1:C:181:ALA:O	1:C:185:VAL:HG23	2.07	0.54
1:B:89:LYS:O	1:B:157:GLN:HG3	2.08	0.54
1:A:205:SER:OG	2:A:3600:PG9:N	2.41	0.53
1:A:548:ILE:HB	1:A:613:MET:HB3	1.90	0.53
1:C:204:SER:HA	1:C:228:GLU:O	2.09	0.53
1:C:438:CYS:CB	1:C:442:CYS:HG	2.21	0.53
1:D:221:ALA:O	1:D:223:LYS:HG2	2.09	0.53
1:A:232:VAL:HG11	1:A:318:ASP:HB3	1.92	0.52
1:C:438:CYS:HG	1:C:442:CYS:HG	1.57	0.52
1:D:535:ALA:HB2	1:D:592:THR:HA	1.92	0.52
1:B:535:ALA:HB2	1:B:592:THR:HA	1.92	0.52
1:B:200:GLY:HA3	1:B:406:LEU:HD11	1.91	0.51
1:D:480:PHE:HE1	1:D:622:PRO:HD2	1.74	0.51
1:B:78:ARG:HB3	1:B:94:ILE:HB	1.92	0.51
1:A:161:VAL:HB	1:A:164:ARG:HG3	1.92	0.51
1:C:88:VAL:HG13	1:C:157:GLN:HB2	1.93	0.51
1:C:373:VAL:O	1:C:394:ARG:NH1	2.43	0.51
1:D:416:PRO:HB2	1:D:431:TYR:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:457:LEU:HB2	1:C:515:THR:HG22	1.91	0.51
1:D:164:ARG:NH2	1:D:175:THR:O	2.43	0.51
1:C:477:PRO:HB2	1:C:621:PHE:HB2	1.91	0.51
1:D:621:PHE:CG	1:D:622:PRO:HA	2.46	0.51
1:A:340:GLU:HB2	1:A:370:HIS:CG	2.46	0.51
1:D:205:SER:HA	1:D:229:SER:O	2.11	0.51
1:A:205:SER:HG	2:A:3600:PG9:H2	1.59	0.51
1:C:438:CYS:HB3	1:C:442:CYS:SG	2.50	0.51
1:C:229:SER:OG	1:C:338:ASP:OD1	2.29	0.51
1:B:274:ASP:HB2	1:B:486:ALA:HA	1.93	0.51
1:C:621:PHE:CD1	1:C:622:PRO:HA	2.46	0.50
1:A:481:ILE:HD11	1:A:485:PHE:CD1	2.46	0.50
1:D:340:GLU:HB2	1:D:370:HIS:CG	2.46	0.50
1:C:404:GLU:HG3	1:C:412:SER:HA	1.93	0.50
1:D:340:GLU:OE2	2:D:3598:PG9:HA	2.12	0.50
1:B:340:GLU:OE2	2:B:3599:PG9:HA	2.11	0.50
1:A:301:ARG:HH11	1:A:301:ARG:HG3	1.76	0.50
1:B:508:PRO:HG2	1:C:508:PRO:HG2	1.94	0.50
1:B:574:ARG:HG2	1:B:576:ARG:HG2	1.94	0.50
1:C:340:GLU:HB2	1:C:370:HIS:CG	2.46	0.49
1:A:481:ILE:HG21	1:A:496:TRP:HB2	1.94	0.49
1:D:70:ASP:HA	1:D:73:ARG:HD3	1.94	0.49
1:B:308:TYR:O	1:B:313:GLN:NE2	2.45	0.49
1:B:111:PRO:HG3	1:B:206:TYR:CG	2.47	0.49
1:A:104:PRO:HB3	1:A:406:LEU:HB3	1.94	0.49
1:A:83:PRO:HA	1:A:89:LYS:HA	1.94	0.49
1:B:457:LEU:HB2	1:B:515:THR:CG2	2.42	0.49
1:C:111:PRO:HG3	1:C:206:TYR:CG	2.48	0.49
1:C:449:LEU:HD21	1:C:457:LEU:HD13	1.95	0.49
1:C:232:VAL:HB	1:C:344:GLY:HA2	1.94	0.49
1:A:328:VAL:O	1:A:330:MET:HG3	2.13	0.49
1:A:92:THR:HG23	1:A:148:GLN:HB2	1.94	0.49
1:D:434:TRP:CG	1:D:435:PRO:HA	2.47	0.49
1:C:274:ASP:HB2	1:C:486:ALA:HA	1.95	0.49
1:C:107:LEU:HD12	1:C:108:THR:N	2.28	0.49
1:D:651:HIS:CD2	1:D:656:GLU:HB2	2.48	0.48
1:D:246:PHE:O	1:D:624:TYR:HA	2.12	0.48
1:A:209:PHE:HB2	1:A:231:MET:HB2	1.94	0.48
1:D:112:TYR:OH	2:D:3598:PG9:OXT	2.20	0.48
1:D:88:VAL:HG13	1:D:157:GLN:HB2	1.94	0.48
1:C:85:ARG:CZ	1:C:179:THR:HB	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:604:HIS:CG	1:D:605:VAL:H	2.32	0.48
1:B:621:PHE:CD1	1:B:622:PRO:HA	2.49	0.48
1:A:628:PRO:HA	1:A:643:TYR:CD1	2.49	0.48
1:C:416:PRO:HB2	1:C:431:TYR:O	2.14	0.48
1:A:91:TYR:HE2	1:A:93:VAL:CG2	2.26	0.48
1:A:110:THR:HB	1:A:207:GLU:HG2	1.96	0.47
1:A:621:PHE:CD1	1:A:622:PRO:HA	2.50	0.47
1:A:532:LEU:HA	1:A:658:THR:HG23	1.96	0.47
1:C:242:HIS:CD2	1:C:247:ARG:HD2	2.48	0.47
1:C:469:SER:HA	1:C:648:GLN:O	2.13	0.47
1:D:438:CYS:HG	1:D:442:CYS:CB	2.27	0.47
1:B:532:LEU:HA	1:B:658:THR:HG23	1.95	0.47
1:A:67:MET:HA	1:A:68:PRO:HD3	1.78	0.47
1:D:479:PRO:HA	1:D:621:PHE:O	2.14	0.47
1:C:200:GLY:HA3	1:C:406:LEU:HD11	1.95	0.47
1:C:549:ASP:OD1	1:C:604:HIS:NE2	2.45	0.47
1:C:81:MET:HB3	1:C:89:LYS:HB3	1.97	0.47
1:A:111:PRO:HG3	1:A:206:TYR:CD2	2.49	0.47
1:B:621:PHE:CG	1:B:622:PRO:HA	2.50	0.47
1:B:546:LYS:HB2	1:B:615:GLN:HB2	1.96	0.47
1:D:94:ILE:HG23	1:D:146:VAL:HG22	1.97	0.47
1:B:255:VAL:HG12	1:B:265:ASN:O	2.15	0.47
1:B:67:MET:HA	1:B:68:PRO:HD3	1.75	0.47
1:D:629:GLN:NE2	1:D:644:THR:O	2.41	0.47
1:D:111:PRO:HG3	1:D:206:TYR:CG	2.50	0.47
1:D:438:CYS:SG	1:D:442:CYS:CB	3.03	0.47
1:A:480:PHE:CE1	1:A:622:PRO:HD2	2.50	0.47
1:D:477:PRO:HB2	1:D:621:PHE:HB2	1.96	0.47
1:B:479:PRO:HA	1:B:621:PHE:O	2.14	0.46
1:C:331:LEU:HD23	1:C:331:LEU:HA	1.79	0.46
1:B:62:PRO:HB2	1:B:64:SER:O	2.15	0.46
1:A:438:CYS:SG	1:A:442:CYS:HB2	2.55	0.46
1:A:507:ARG:HB3	1:A:509:ASP:OD2	2.16	0.46
1:C:67:MET:HA	1:C:68:PRO:HD3	1.84	0.46
1:C:221:ALA:O	1:C:223:LYS:HG2	2.15	0.46
1:C:556:PRO:HG3	1:D:129:GLU:HA	1.97	0.46
1:B:340:GLU:HB2	1:B:370:HIS:CD2	2.51	0.46
1:D:369:ARG:H	1:D:369:ARG:HG3	1.61	0.46
1:C:120:ARG:CZ	1:D:556:PRO:HB2	2.46	0.46
1:C:480:PHE:CE1	1:C:622:PRO:HD2	2.51	0.46
1:D:480:PHE:CE1	1:D:622:PRO:HD2	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:496:TRP:HA	1:B:499:GLN:HG2	1.96	0.46
1:B:416:PRO:HB2	1:B:431:TYR:O	2.14	0.46
1:C:146:VAL:HG11	1:C:188:LEU:HD11	1.97	0.46
1:C:621:PHE:CG	1:C:622:PRO:HA	2.51	0.46
1:D:85:ARG:N	1:D:183:ASP:OD2	2.48	0.46
1:A:223:LYS:HE3	3:A:2433:GOL:H2	1.97	0.46
1:A:540:ASP:O	1:A:629:GLN:HA	2.16	0.46
1:B:81:MET:HB3	1:B:89:LYS:HB3	1.98	0.46
1:A:574:ARG:HB3	1:A:577:TYR:HD1	1.81	0.46
1:B:205:SER:HB3	1:B:206:TYR:H	1.47	0.46
1:C:481:ILE:HD11	1:C:485:PHE:CD1	2.51	0.46
1:A:274:ASP:HB2	1:A:486:ALA:HA	1.99	0.45
1:C:224:VAL:HG23	1:C:329:PRO:HB2	1.97	0.45
1:B:153:LYS:HD2	1:B:153:LYS:HA	1.66	0.45
1:A:237:GLY:HA2	1:A:242:HIS:HA	1.99	0.45
1:C:334:GLN:HB3	1:C:345:ALA:HB3	1.99	0.45
1:A:480:PHE:HE1	1:A:622:PRO:HD2	1.82	0.45
1:D:91:TYR:CE2	1:D:93:VAL:CG2	2.99	0.45
1:D:70:ASP:OD1	1:D:70:ASP:N	2.49	0.45
1:C:296:TYR:O	1:C:300:GLN:HG3	2.16	0.45
1:C:552:PRO:HD3	1:C:610:HIS:CE1	2.51	0.45
1:B:331:LEU:HD23	1:B:331:LEU:HA	1.69	0.45
1:D:566:LEU:HD12	1:D:567:PRO:HD2	1.99	0.45
1:C:537:SER:HB3	1:C:647:THR:H	1.80	0.45
1:C:73:ARG:HH11	1:C:123:ASN:HB3	1.82	0.45
1:D:478:VAL:HG22	1:D:507:ARG:HE	1.82	0.45
1:A:508:PRO:HG2	1:D:508:PRO:HG2	1.98	0.45
1:B:628:PRO:C	1:B:630:LYS:H	2.18	0.45
1:A:128:ARG:NH2	1:B:557:ASP:OD2	2.29	0.45
1:C:120:ARG:HB3	1:C:129:GLU:O	2.17	0.45
1:B:544:VAL:O	1:B:616:ILE:HA	2.17	0.45
1:B:185:VAL:O	1:B:189:VAL:HG13	2.17	0.45
1:D:511:VAL:HG12	1:D:616:ILE:HB	1.99	0.45
1:B:544:VAL:HB	1:B:617:GLN:HG2	1.99	0.44
1:B:216:LEU:HD23	1:B:321:LEU:HD12	1.99	0.44
1:D:481:ILE:HG21	1:D:496:TRP:HB2	1.99	0.44
1:D:124:ALA:CB	1:D:129:GLU:HB3	2.48	0.44
1:D:67:MET:HA	1:D:68:PRO:HD3	1.79	0.44
1:A:566:LEU:HA	1:A:567:PRO:HD3	1.78	0.44
1:A:271:ASP:HA	1:D:553:ALA:HB1	1.99	0.44
1:A:604:HIS:CG	1:A:605:VAL:H	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:91:TYR:HE2	1:D:93:VAL:HG23	1.83	0.44
1:B:494:LYS:HB2	1:B:495:PRO:HD3	2.00	0.44
1:B:514:GLU:HG3	1:B:611:ARG:NH2	2.33	0.44
1:B:481:ILE:HG21	1:B:496:TRP:HB2	1.99	0.44
1:B:529:VAL:HG11	1:B:596:HIS:CE1	2.53	0.44
1:D:515:THR:O	1:D:611:ARG:NH2	2.51	0.44
1:C:510:VAL:HG22	1:C:617:GLN:HB3	2.00	0.44
1:D:370:HIS:C	1:D:372:GLY:H	2.22	0.43
1:C:106:LEU:HD23	1:C:145:ARG:CZ	2.48	0.43
1:A:121:VAL:HB	1:A:124:ALA:HB2	2.00	0.43
1:B:417:ASP:OD2	1:B:433:SER:HA	2.18	0.43
1:A:146:VAL:HG11	1:A:188:LEU:HD11	2.00	0.43
1:A:224:VAL:HG23	1:A:329:PRO:HB2	2.00	0.43
1:C:177:GLU:HB2	1:C:213:MET:HE2	2.00	0.43
1:A:604:HIS:CG	1:A:605:VAL:N	2.87	0.43
1:B:529:VAL:HB	1:B:661:LEU:HB2	1.99	0.43
1:B:164:ARG:NH2	1:B:173:THR:OG1	2.51	0.43
1:B:523:ARG:HH11	1:B:523:ARG:HG3	1.83	0.43
1:C:574:ARG:HB3	1:C:577:TYR:HD1	1.82	0.43
1:A:334:GLN:HB3	1:A:345:ALA:HB3	2.00	0.43
1:D:205:SER:HB2	1:D:370:HIS:NE2	2.33	0.43
1:C:438:CYS:SG	1:C:442:CYS:SG	3.14	0.43
1:B:502:ARG:HG2	1:B:565:GLU:OE2	2.18	0.43
1:A:535:ALA:HB2	1:A:592:THR:HA	2.00	0.43
1:A:479:PRO:HA	1:A:621:PHE:O	2.19	0.43
1:D:417:ASP:OD2	1:D:433:SER:HA	2.19	0.43
1:C:111:PRO:HB3	1:C:150:ILE:HD11	2.01	0.43
1:A:91:TYR:HB3	1:A:149:ASP:HB2	2.00	0.43
1:D:91:TYR:CE2	1:D:93:VAL:HG23	2.53	0.43
1:B:406:LEU:O	1:B:408:PRO:HD3	2.19	0.43
1:D:117:ARG:HA	1:D:117:ARG:HD2	1.89	0.43
1:D:110:THR:HB	1:D:207:GLU:HG2	2.01	0.43
1:B:480:PHE:HD2	1:B:496:TRP:CE3	2.37	0.42
1:A:521:PRO:HA	1:A:606:PHE:O	2.19	0.42
1:C:318:ASP:OD2	1:C:319:LYS:N	2.52	0.42
1:C:527:VAL:HA	1:C:528:PRO:HD3	1.90	0.42
1:B:566:LEU:HD12	1:B:567:PRO:HD2	2.01	0.42
1:D:153:LYS:HD2	1:D:261:ARG:O	2.19	0.42
1:D:662:LEU:HA	1:D:663:PRO:HD3	1.82	0.42
1:D:205:SER:OG	2:D:3598:PG9:C	2.68	0.42
1:A:221:ALA:O	1:A:223:LYS:HG2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:511:VAL:HG11	1:B:650:ILE:HG13	2.02	0.42
1:C:425:ASP:OD2	1:C:523:ARG:NH2	2.39	0.42
1:B:497:LEU:HD23	1:B:497:LEU:HA	1.75	0.42
1:A:365:MET:HB2	1:A:420:ILE:HG23	2.01	0.42
1:C:232:VAL:HG13	1:C:317:LEU:HB2	2.01	0.42
1:B:558:ASP:OD1	1:B:560:LYS:HD3	2.19	0.42
1:A:540:ASP:OD2	1:A:576:ARG:NH2	2.46	0.42
1:D:494:LYS:HB2	1:D:495:PRO:HD3	2.02	0.42
1:D:228:GLU:O	1:D:229:SER:C	2.58	0.42
1:A:153:LYS:HA	1:A:153:LYS:HD2	1.90	0.42
1:D:442:CYS:O	1:D:443:THR:C	2.58	0.41
1:C:185:VAL:HG12	1:C:221:ALA:HB3	2.02	0.41
1:A:339:GLN:HE22	1:A:370:HIS:H	1.66	0.41
1:D:205:SER:OG	2:D:3598:PG9:OXT	2.38	0.41
1:C:107:LEU:HD12	1:C:108:THR:H	1.85	0.41
1:C:343:TRP:CD2	1:C:347:HIS:CE1	3.08	0.41
1:A:502:ARG:NH2	1:D:484:PRO:HG3	2.35	0.41
1:C:583:LYS:HA	1:C:584:PRO:HD2	1.92	0.41
1:C:153:LYS:HD2	1:C:261:ARG:O	2.21	0.41
1:B:477:PRO:HB2	1:B:621:PHE:HB2	2.02	0.41
1:A:55:VAL:HG22	1:C:170:LEU:HA	2.01	0.41
1:D:453:ASP:N	1:D:453:ASP:OD1	2.54	0.41
1:A:82:VAL:O	1:A:90:LEU:N	2.40	0.41
1:A:315:GLN:O	1:A:317:LEU:N	2.54	0.41
1:B:221:ALA:O	1:B:223:LYS:HG2	2.19	0.41
1:B:91:TYR:HB3	1:B:149:ASP:HB2	2.02	0.41
1:C:422:ASN:HB2	1:C:429:ASP:OD2	2.20	0.41
1:D:267:ILE:HA	1:D:268:PRO:HD2	1.89	0.41
1:D:523:ARG:HG2	1:D:605:VAL:HG22	2.01	0.41
1:D:324:ARG:HA	1:D:324:ARG:HD3	1.50	0.41
1:C:551:GLN:HG3	1:C:564:TYR:HB3	2.03	0.41
1:B:401:PHE:HA	1:B:413:VAL:HG21	2.01	0.41
1:B:85:ARG:CZ	1:B:179:THR:HB	2.51	0.41
1:B:164:ARG:HA	1:B:165:PRO:HD3	1.88	0.41
1:C:369:ARG:HG3	1:C:369:ARG:H	1.59	0.41
1:A:512:THR:HG23	1:A:615:GLN:HG3	2.02	0.41
1:D:574:ARG:HB3	1:D:577:TYR:CD1	2.55	0.41
1:B:480:PHE:HE1	1:B:622:PRO:HD2	1.85	0.41
1:B:545:VAL:HG11	1:B:660:ILE:HD11	2.03	0.41
1:A:434:TRP:CG	1:A:435:PRO:HA	2.55	0.41
1:C:121:VAL:HA	1:C:122:PRO:HD2	1.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:LEU:HD23	1:A:296:TYR:CD2	2.56	0.41
1:D:73:ARG:HG3	1:D:125:LEU:HA	2.03	0.40
1:C:529:VAL:HB	1:C:661:LEU:HB2	2.03	0.40
1:B:124:ALA:CB	1:B:129:GLU:HB3	2.51	0.40
1:B:209:PHE:HB2	1:B:231:MET:HB2	2.04	0.40
1:C:621:PHE:HA	1:C:622:PRO:HA	1.83	0.40
1:A:136:ASP:OD1	1:A:136:ASP:N	2.54	0.40
1:D:178:THR:HG23	1:D:213:MET:O	2.21	0.40
1:C:228:GLU:O	1:C:229:SER:C	2.57	0.40
1:A:191:ASN:O	1:A:193:PRO:HD3	2.21	0.40
1:C:53:LEU:HD23	1:C:53:LEU:HA	1.88	0.40
1:B:324:ARG:HA	1:B:324:ARG:HD3	1.85	0.40
1:B:280:LYS:HB2	1:B:280:LYS:HE2	1.48	0.40
1:D:451:LEU:HA	1:D:457:LEU:HD23	2.03	0.40
1:B:228:GLU:O	1:B:229:SER:C	2.59	0.40
1:B:399:ARG:HB3	1:B:400:PRO:HD3	2.04	0.40
1:B:583:LYS:HA	1:B:584:PRO:HD2	1.93	0.40
1:B:232:VAL:HG11	1:B:318:ASP:HB3	2.04	0.40
1:D:331:LEU:HD23	1:D:362:THR:HB	2.04	0.40
1:D:540:ASP:OD2	1:D:576:ARG:NH1	2.52	0.40
1:A:583:LYS:HA	1:A:584:PRO:HD2	1.93	0.40
1:A:162:MET:HB3	1:A:298:PHE:CD1	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	615/652 (94%)	558 (91%)	53 (9%)	4 (1%)	26 66
1	B	615/652 (94%)	568 (92%)	46 (8%)	1 (0%)	52 85
1	C	615/652 (94%)	569 (92%)	41 (7%)	5 (1%)	24 62

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	615/652 (94%)	558 (91%)	53 (9%)	4 (1%)	26	66
All	All	2460/2608 (94%)	2253 (92%)	193 (8%)	14 (1%)	30	68

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	344	GLY
1	A	316	ALA
1	A	344	GLY
1	B	344	GLY
1	C	344	GLY
1	C	316	ALA
1	A	237	GLY
1	C	441	ASN
1	D	345	ALA
1	D	371	SER
1	A	495	PRO
1	C	113	ASN
1	D	495	PRO
1	C	237	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	517/544 (95%)	470 (91%)	47 (9%)	12	41
1	B	517/544 (95%)	471 (91%)	46 (9%)	12	43
1	C	517/544 (95%)	464 (90%)	53 (10%)	9	34
1	D	517/544 (95%)	461 (89%)	56 (11%)	8	32
All	All	2068/2176 (95%)	1866 (90%)	202 (10%)	10	37

All (202) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	SER
1	A	56	GLN
1	A	64	SER
1	A	65	VAL
1	A	67	MET
1	A	70	ASP
1	A	71	GLN
1	A	77	LYS
1	A	110	THR
1	A	120	ARG
1	A	130	VAL
1	A	149	ASP
1	A	174	LYS
1	A	195	SER
1	A	198	ARG
1	A	205	SER
1	A	229	SER
1	A	239	ASP
1	A	256	SER
1	A	258	MET
1	A	319	LYS
1	A	339	GLN
1	A	398	PHE
1	A	399	ARG
1	A	427	LYS
1	A	432	ARG
1	A	440	SER
1	A	442	CYS
1	A	447	THR
1	A	453	ASP
1	A	472	SER
1	A	490	SER
1	A	491	SER
1	A	511	VAL
1	A	523	ARG
1	A	524	VAL
1	A	536	THR
1	A	539	THR
1	A	541	SER
1	A	576	ARG
1	A	615	GLN
1	A	620	TRP
1	A	621	PHE

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Mol	Chain	Res	Type
1	A	629	GLN
1	A	630	LYS
1	A	632	VAL
1	A	644	THR
1	B	57	THR
1	B	59	SER
1	B	64	SER
1	B	65	VAL
1	B	67	MET
1	B	70	ASP
1	B	77	LYS
1	B	80	VAL
1	B	99	ASN
1	B	107	LEU
1	B	120	ARG
1	B	126	THR
1	B	157	GLN
1	B	201	MET
1	B	205	SER
1	B	256	SER
1	B	276	THR
1	B	279	LEU
1	B	280	LYS
1	B	308	TYR
1	B	313	GLN
1	B	319	LYS
1	B	325	LYS
1	B	331	LEU
1	B	379	THR
1	B	395	ARG
1	B	412	SER
1	B	415	LEU
1	B	425	ASP
1	B	438	CYS
1	B	443	THR
1	B	469	SER
1	B	472	SER
1	B	482	SER
1	B	519	ASP
1	B	520	GLU
1	B	539	THR
1	B	560	LYS

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Mol	Chain	Res	Type
1	B	565	GLU
1	B	598	THR
1	B	620	TRP
1	B	629	GLN
1	B	648	GLN
1	B	649	SER
1	B	665	VAL
1	B	666	LYS
1	C	54	SER
1	C	57	THR
1	C	64	SER
1	C	65	VAL
1	C	66	HIS
1	C	71	GLN
1	C	77	LYS
1	C	85	ARG
1	C	98	LYS
1	C	120	ARG
1	C	123	ASN
1	C	129	GLU
1	C	130	VAL
1	C	140	GLU
1	C	174	LYS
1	C	175	THR
1	C	198	ARG
1	C	205	SER
1	C	206	TYR
1	C	209	PHE
1	C	270	ARG
1	C	330	MET
1	C	331	LEU
1	C	346	ILE
1	C	357	VAL
1	C	379	THR
1	C	398	PHE
1	C	412	SER
1	C	427	LYS
1	C	432	ARG
1	C	438	CYS
1	C	453	ASP
1	C	465	ASP
1	C	482	SER

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Mol	Chain	Res	Type
1	C	490	SER
1	C	491	SER
1	C	506	SER
1	C	518	LEU
1	C	523	ARG
1	C	537	SER
1	C	541	SER
1	C	560	LYS
1	C	576	ARG
1	C	583	LYS
1	C	615	GLN
1	C	616	ILE
1	C	620	TRP
1	C	621	PHE
1	C	629	GLN
1	C	644	THR
1	C	648	GLN
1	C	649	SER
1	C	655	LYS
1	D	61	ILE
1	D	64	SER
1	D	65	VAL
1	D	69	THR
1	D	70	ASP
1	D	71	GLN
1	D	72	GLN
1	D	73	ARG
1	D	77	LYS
1	D	80	VAL
1	D	101	ARG
1	D	120	ARG
1	D	130	VAL
1	D	153	LYS
1	D	204	SER
1	D	205	SER
1	D	206	TYR
1	D	265	ASN
1	D	266	ASP
1	D	280	LYS
1	D	293	LEU
1	D	308	TYR
1	D	324	ARG

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Mol	Chain	Res	Type
1	D	325	LYS
1	D	327	THR
1	D	331	LEU
1	D	358	LYS
1	D	365	MET
1	D	379	THR
1	D	395	ARG
1	D	398	PHE
1	D	417	ASP
1	D	436	SER
1	D	440	SER
1	D	442	CYS
1	D	447	THR
1	D	449	LEU
1	D	472	SER
1	D	482	SER
1	D	491	SER
1	D	506	SER
1	D	518	LEU
1	D	519	ASP
1	D	539	THR
1	D	541	SER
1	D	576	ARG
1	D	611	ARG
1	D	620	TRP
1	D	621	PHE
1	D	626	ARG
1	D	630	LYS
1	D	635	ILE
1	D	644	THR
1	D	649	SER
1	D	655	LYS
1	D	665	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	A	133	GLN
1	A	191	ASN
1	A	339	GLN
1	B	72	GLN

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Mol	Chain	Res	Type
1	B	133	GLN
1	B	392	GLN
1	B	455	HIS
1	B	629	GLN
1	B	651	HIS
1	C	71	GLN
1	C	190	HIS
1	D	71	GLN
1	D	102	ASN
1	D	133	GLN
1	D	594	HIS
1	D	651	HIS

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](#)

7 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	GOL	A	2431	-	5,5,5	0.28	0	5,5,5	0.26	0
3	GOL	A	2433	-	5,5,5	0.34	0	5,5,5	0.26	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PG9	A	3600	-	7,11,11	1.73	1 (14%)	8,14,14	0.96	0
3	GOL	B	2432	-	5,5,5	0.26	0	5,5,5	0.27	0
2	PG9	B	3599	-	7,11,11	2.12	1 (14%)	8,14,14	1.02	0
3	GOL	C	2430	-	5,5,5	0.36	0	5,5,5	0.16	0
2	PG9	D	3598	-	7,11,11	1.93	1 (14%)	8,14,14	0.98	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	GOL	A	2431	-	-	0/4/4/4	0/0/0/0
3	GOL	A	2433	-	-	0/4/4/4	0/0/0/0
2	PG9	A	3600	-	-	0/4/8/8	0/1/1/1
3	GOL	B	2432	-	-	0/4/4/4	0/0/0/0
2	PG9	B	3599	-	-	0/4/8/8	0/1/1/1
3	GOL	C	2430	-	-	0/4/4/4	0/0/0/0
2	PG9	D	3598	-	-	0/4/8/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3599	PG9	CB-CA	-5.48	1.47	1.52
2	D	3598	PG9	CB-CA	-4.94	1.47	1.52
2	A	3600	PG9	CB-CA	-4.29	1.48	1.52

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2433	GOL	1	0
2	A	3600	PG9	2	0
2	B	3599	PG9	1	0
2	D	3598	PG9	4	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	617/652 (94%)	-0.42	6 (0%) 84 80	19, 19, 19, 19	0
1	B	617/652 (94%)	-0.40	7 (1%) 82 78	19, 19, 19, 19	0
1	C	617/652 (94%)	-0.40	11 (1%) 71 65	19, 19, 19, 19	0
1	D	617/652 (94%)	-0.40	12 (1%) 70 63	19, 19, 19, 19	0
All	All	2468/2608 (94%)	-0.41	36 (1%) 76 71	19, 19, 19, 19	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	70	ASP	4.1
1	A	66	HIS	4.0
1	A	50	HIS	3.7
1	D	66	HIS	3.4
1	D	50	HIS	3.3
1	B	99	ASN	3.1
1	C	66	HIS	3.1
1	B	50	HIS	3.1
1	B	69	THR	3.1
1	A	70	ASP	3.0
1	B	441	ASN	2.9
1	D	99	ASN	2.9
1	D	70	ASP	2.9
1	C	70	ASP	2.8
1	C	441	ASN	2.8
1	B	66	HIS	2.7
1	A	69	THR	2.7
1	D	69	THR	2.7
1	C	64	SER	2.6
1	A	68	PRO	2.6
1	D	441	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	50	HIS	2.5
1	D	71	GLN	2.5
1	C	101	ARG	2.4
1	D	440	SER	2.4
1	D	63	ALA	2.3
1	C	99	ASN	2.3
1	C	63	ALA	2.3
1	A	64	SER	2.2
1	D	101	ARG	2.2
1	C	414	HIS	2.2
1	C	72	GLN	2.1
1	D	409	GLY	2.1
1	D	174	LYS	2.1
1	B	265	ASN	2.1
1	C	519	ASP	2.1

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	PG9	B	3599	11/11	0.92	0.42	12.57	19,19,19,19	0
2	PG9	A	3600	11/11	0.90	0.42	8.61	19,19,19,19	0
2	PG9	D	3598	11/11	0.95	0.33	6.41	19,19,19,19	0
3	GOL	B	2432	6/6	0.92	0.24	3.41	19,19,19,19	0
3	GOL	C	2430	6/6	0.91	0.15	-	19,19,19,19	0
3	GOL	A	2431	6/6	0.91	0.19	-	19,19,19,19	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	GOL	A	2433	6/6	0.81	0.40	-	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.