



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 02:16 AM UTC

PDB ID : 5Z19 / pdb_00005z19
Title : The crystal structure of Ruminococcus gnavus beta-glucuronidase in complex with uronic isofagomine
Authors : Dashnyam, P.; Lin, H.Y.; Lin, C.H.
Deposited on : 2017-12-25
Resolution : 2.50 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

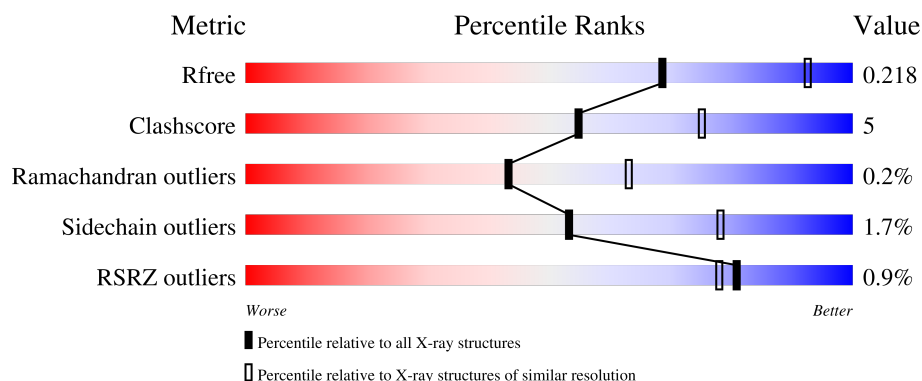
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	627	% 83% 11% 5%
1	B	627	% 82% 12% • •
1	C	627	% 86% 10% •
1	D	627	% 82% 12% • 5%
1	E	627	% 84% 11% • 5%

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Mol	Chain	Length	Quality of chain
1	F	627	% 83% 11% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 29830 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 Beta-glucuronidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	593	Total	C	N	O	S	0	0	0
			4864	3138	791	911	24			
1	B	601	Total	C	N	O	S	0	0	0
			4923	3173	803	923	24			
1	C	600	Total	C	N	O	S	0	0	0
			4915	3167	802	922	24			
1	D	596	Total	C	N	O	S	0	0	0
			4880	3146	795	915	24			
1	E	597	Total	C	N	O	S	0	0	0
			4889	3152	797	916	24			
1	F	600	Total	C	N	O	S	0	0	0
			4915	3167	802	922	24			

There are 150 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	initiating methionine	UNP Q6W7J7
A	-22	HIS	-	expression tag	UNP Q6W7J7
A	-21	HIS	-	expression tag	UNP Q6W7J7
A	-20	HIS	-	expression tag	UNP Q6W7J7
A	-19	HIS	-	expression tag	UNP Q6W7J7
A	-18	HIS	-	expression tag	UNP Q6W7J7
A	-17	HIS	-	expression tag	UNP Q6W7J7
A	-16	SER	-	expression tag	UNP Q6W7J7
A	-15	SER	-	expression tag	UNP Q6W7J7
A	-14	GLY	-	expression tag	UNP Q6W7J7
A	-13	VAL	-	expression tag	UNP Q6W7J7
A	-12	ASP	-	expression tag	UNP Q6W7J7
A	-11	LEU	-	expression tag	UNP Q6W7J7
A	-10	GLY	-	expression tag	UNP Q6W7J7
A	-9	THR	-	expression tag	UNP Q6W7J7
A	-8	GLU	-	expression tag	UNP Q6W7J7
A	-7	ASN	-	expression tag	UNP Q6W7J7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	LEU	-	expression tag	UNP Q6W7J7
A	-5	TYR	-	expression tag	UNP Q6W7J7
A	-4	PHE	-	expression tag	UNP Q6W7J7
A	-3	GLN	-	expression tag	UNP Q6W7J7
A	-2	SER	-	expression tag	UNP Q6W7J7
A	-1	ASN	-	expression tag	UNP Q6W7J7
A	0	GLY	-	expression tag	UNP Q6W7J7
A	1	MET	-	expression tag	UNP Q6W7J7
B	-23	MET	-	initiating methionine	UNP Q6W7J7
B	-22	HIS	-	expression tag	UNP Q6W7J7
B	-21	HIS	-	expression tag	UNP Q6W7J7
B	-20	HIS	-	expression tag	UNP Q6W7J7
B	-19	HIS	-	expression tag	UNP Q6W7J7
B	-18	HIS	-	expression tag	UNP Q6W7J7
B	-17	HIS	-	expression tag	UNP Q6W7J7
B	-16	SER	-	expression tag	UNP Q6W7J7
B	-15	SER	-	expression tag	UNP Q6W7J7
B	-14	GLY	-	expression tag	UNP Q6W7J7
B	-13	VAL	-	expression tag	UNP Q6W7J7
B	-12	ASP	-	expression tag	UNP Q6W7J7
B	-11	LEU	-	expression tag	UNP Q6W7J7
B	-10	GLY	-	expression tag	UNP Q6W7J7
B	-9	THR	-	expression tag	UNP Q6W7J7
B	-8	GLU	-	expression tag	UNP Q6W7J7
B	-7	ASN	-	expression tag	UNP Q6W7J7
B	-6	LEU	-	expression tag	UNP Q6W7J7
B	-5	TYR	-	expression tag	UNP Q6W7J7
B	-4	PHE	-	expression tag	UNP Q6W7J7
B	-3	GLN	-	expression tag	UNP Q6W7J7
B	-2	SER	-	expression tag	UNP Q6W7J7
B	-1	ASN	-	expression tag	UNP Q6W7J7
B	0	GLY	-	expression tag	UNP Q6W7J7
B	1	MET	-	expression tag	UNP Q6W7J7
C	-23	MET	-	initiating methionine	UNP Q6W7J7
C	-22	HIS	-	expression tag	UNP Q6W7J7
C	-21	HIS	-	expression tag	UNP Q6W7J7
C	-20	HIS	-	expression tag	UNP Q6W7J7
C	-19	HIS	-	expression tag	UNP Q6W7J7
C	-18	HIS	-	expression tag	UNP Q6W7J7
C	-17	HIS	-	expression tag	UNP Q6W7J7
C	-16	SER	-	expression tag	UNP Q6W7J7
C	-15	SER	-	expression tag	UNP Q6W7J7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-14	GLY	-	expression tag	UNP Q6W7J7
C	-13	VAL	-	expression tag	UNP Q6W7J7
C	-12	ASP	-	expression tag	UNP Q6W7J7
C	-11	LEU	-	expression tag	UNP Q6W7J7
C	-10	GLY	-	expression tag	UNP Q6W7J7
C	-9	THR	-	expression tag	UNP Q6W7J7
C	-8	GLU	-	expression tag	UNP Q6W7J7
C	-7	ASN	-	expression tag	UNP Q6W7J7
C	-6	LEU	-	expression tag	UNP Q6W7J7
C	-5	TYR	-	expression tag	UNP Q6W7J7
C	-4	PHE	-	expression tag	UNP Q6W7J7
C	-3	GLN	-	expression tag	UNP Q6W7J7
C	-2	SER	-	expression tag	UNP Q6W7J7
C	-1	ASN	-	expression tag	UNP Q6W7J7
C	0	GLY	-	expression tag	UNP Q6W7J7
C	1	MET	-	expression tag	UNP Q6W7J7
D	-23	MET	-	initiating methionine	UNP Q6W7J7
D	-22	HIS	-	expression tag	UNP Q6W7J7
D	-21	HIS	-	expression tag	UNP Q6W7J7
D	-20	HIS	-	expression tag	UNP Q6W7J7
D	-19	HIS	-	expression tag	UNP Q6W7J7
D	-18	HIS	-	expression tag	UNP Q6W7J7
D	-17	HIS	-	expression tag	UNP Q6W7J7
D	-16	SER	-	expression tag	UNP Q6W7J7
D	-15	SER	-	expression tag	UNP Q6W7J7
D	-14	GLY	-	expression tag	UNP Q6W7J7
D	-13	VAL	-	expression tag	UNP Q6W7J7
D	-12	ASP	-	expression tag	UNP Q6W7J7
D	-11	LEU	-	expression tag	UNP Q6W7J7
D	-10	GLY	-	expression tag	UNP Q6W7J7
D	-9	THR	-	expression tag	UNP Q6W7J7
D	-8	GLU	-	expression tag	UNP Q6W7J7
D	-7	ASN	-	expression tag	UNP Q6W7J7
D	-6	LEU	-	expression tag	UNP Q6W7J7
D	-5	TYR	-	expression tag	UNP Q6W7J7
D	-4	PHE	-	expression tag	UNP Q6W7J7
D	-3	GLN	-	expression tag	UNP Q6W7J7
D	-2	SER	-	expression tag	UNP Q6W7J7
D	-1	ASN	-	expression tag	UNP Q6W7J7
D	0	GLY	-	expression tag	UNP Q6W7J7
D	1	MET	-	expression tag	UNP Q6W7J7
E	-23	MET	-	initiating methionine	UNP Q6W7J7

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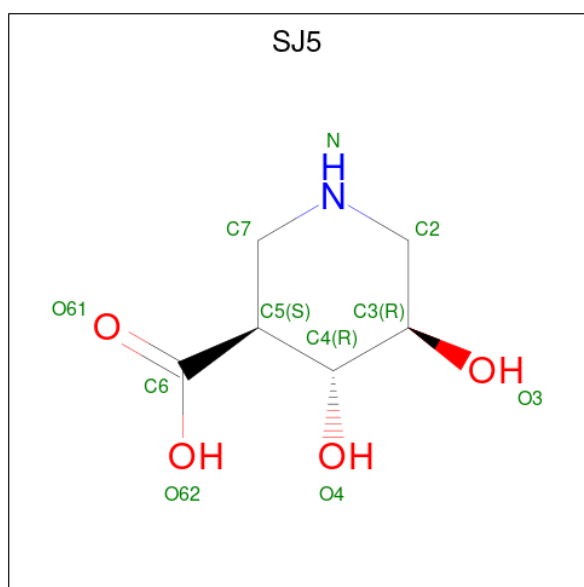
Chain	Residue	Modelled	Actual	Comment	Reference
E	-22	HIS	-	expression tag	UNP Q6W7J7
E	-21	HIS	-	expression tag	UNP Q6W7J7
E	-20	HIS	-	expression tag	UNP Q6W7J7
E	-19	HIS	-	expression tag	UNP Q6W7J7
E	-18	HIS	-	expression tag	UNP Q6W7J7
E	-17	HIS	-	expression tag	UNP Q6W7J7
E	-16	SER	-	expression tag	UNP Q6W7J7
E	-15	SER	-	expression tag	UNP Q6W7J7
E	-14	GLY	-	expression tag	UNP Q6W7J7
E	-13	VAL	-	expression tag	UNP Q6W7J7
E	-12	ASP	-	expression tag	UNP Q6W7J7
E	-11	LEU	-	expression tag	UNP Q6W7J7
E	-10	GLY	-	expression tag	UNP Q6W7J7
E	-9	THR	-	expression tag	UNP Q6W7J7
E	-8	GLU	-	expression tag	UNP Q6W7J7
E	-7	ASN	-	expression tag	UNP Q6W7J7
E	-6	LEU	-	expression tag	UNP Q6W7J7
E	-5	TYR	-	expression tag	UNP Q6W7J7
E	-4	PHE	-	expression tag	UNP Q6W7J7
E	-3	GLN	-	expression tag	UNP Q6W7J7
E	-2	SER	-	expression tag	UNP Q6W7J7
E	-1	ASN	-	expression tag	UNP Q6W7J7
E	0	GLY	-	expression tag	UNP Q6W7J7
E	1	MET	-	expression tag	UNP Q6W7J7
F	-23	MET	-	initiating methionine	UNP Q6W7J7
F	-22	HIS	-	expression tag	UNP Q6W7J7
F	-21	HIS	-	expression tag	UNP Q6W7J7
F	-20	HIS	-	expression tag	UNP Q6W7J7
F	-19	HIS	-	expression tag	UNP Q6W7J7
F	-18	HIS	-	expression tag	UNP Q6W7J7
F	-17	HIS	-	expression tag	UNP Q6W7J7
F	-16	SER	-	expression tag	UNP Q6W7J7
F	-15	SER	-	expression tag	UNP Q6W7J7
F	-14	GLY	-	expression tag	UNP Q6W7J7
F	-13	VAL	-	expression tag	UNP Q6W7J7
F	-12	ASP	-	expression tag	UNP Q6W7J7
F	-11	LEU	-	expression tag	UNP Q6W7J7
F	-10	GLY	-	expression tag	UNP Q6W7J7
F	-9	THR	-	expression tag	UNP Q6W7J7
F	-8	GLU	-	expression tag	UNP Q6W7J7
F	-7	ASN	-	expression tag	UNP Q6W7J7
F	-6	LEU	-	expression tag	UNP Q6W7J7

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-5	TYR	-	expression tag	UNP Q6W7J7
F	-4	PHE	-	expression tag	UNP Q6W7J7
F	-3	GLN	-	expression tag	UNP Q6W7J7
F	-2	SER	-	expression tag	UNP Q6W7J7
F	-1	ASN	-	expression tag	UNP Q6W7J7
F	0	GLY	-	expression tag	UNP Q6W7J7
F	1	MET	-	expression tag	UNP Q6W7J7

- Molecule 2 is (3S,4R,5R)-4,5-dihydroxypiperidine-3-carboxylic acid (CCD ID: SJ5) (formula: $C_6H_{11}NO_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	6	1	4		
2	B	1	Total	C	N	O	0	0
			11	6	1	4		
2	C	1	Total	C	N	O	0	0
			11	6	1	4		
2	D	1	Total	C	N	O	0	0
			11	6	1	4		
2	E	1	Total	C	N	O	0	0
			11	6	1	4		
2	F	1	Total	C	N	O	0	0
			11	6	1	4		

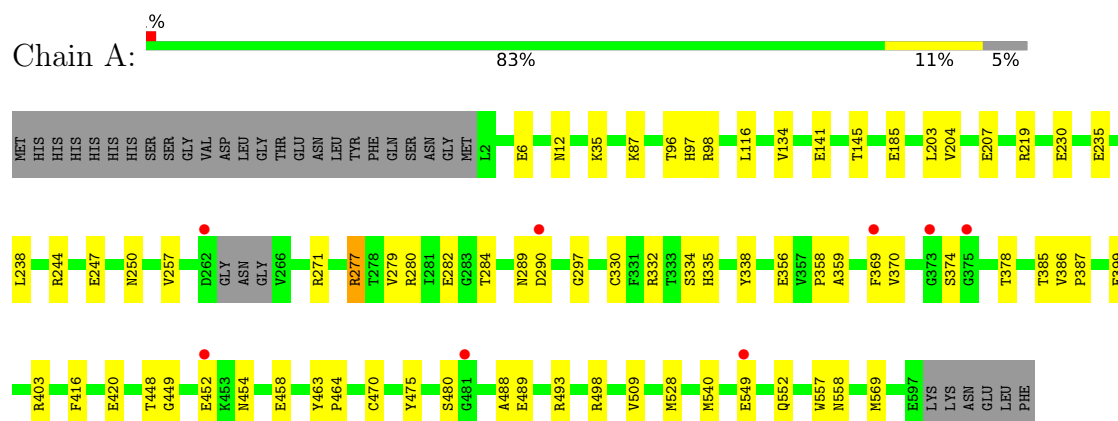
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	58	Total 58	O 58	0	0
3	B	64	Total 64	O 64	0	0
3	C	72	Total 72	O 72	0	0
3	D	59	Total 59	O 59	0	0
3	E	59	Total 59	O 59	0	0
3	F	66	Total 66	O 66	0	0

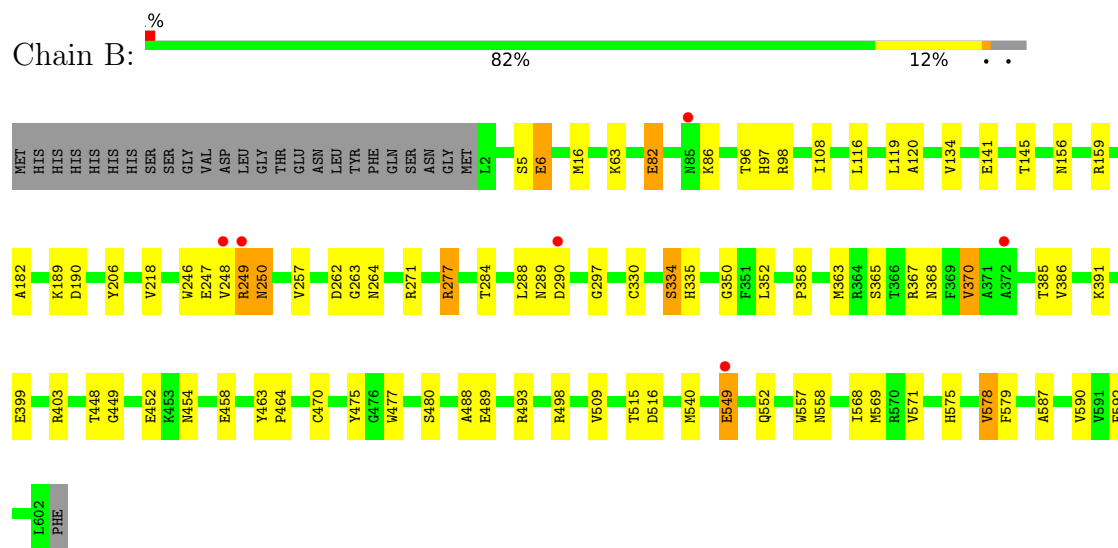
3 Residue-property plots [i](#)

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.

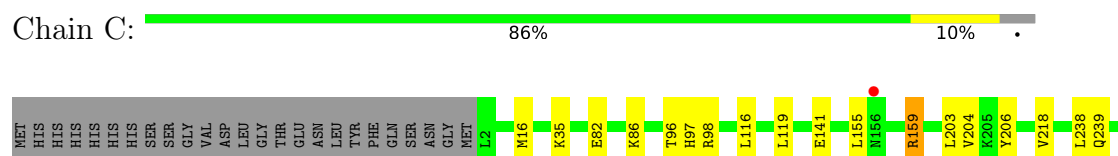
• Molecule 1: Beta-glucuronidase

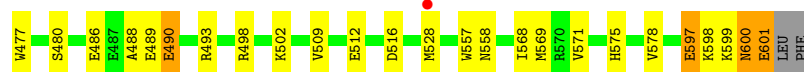


• Molecule 1: Beta-glucuronidase



• Molecule 1: Beta-glucuronidase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	181.28Å 112.36Å 209.79Å 90.00° 93.45° 90.00°	Depositor
Resolution (Å)	29.92 – 2.50 29.92 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.5 (29.92-2.50) 98.4 (29.92-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.166 , 0.209 0.186 , 0.218	Depositor DCC
R_{free} test set	7142 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	44.6	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 26.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	29830	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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: SJ5

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	0.31	0/5003	0.72	0/6784
1	B	0.37	0/5063	0.73	2/6864 (0.0%)
1	C	0.35	0/5055	0.75	1/6853 (0.0%)
1	D	0.38	0/5020	0.75	3/6808 (0.0%)
1	E	0.32	0/5029	0.73	3/6819 (0.0%)
1	F	0.33	0/5055	0.73	1/6853 (0.0%)
All	All	0.35	0/30225	0.74	10/40981 (0.0%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	263	GLY	CA-C-N	6.68	133.73	121.70
1	D	263	GLY	C-N-CA	6.68	133.73	121.70
1	F	4	TYR	N-CA-C	-6.68	98.92	109.07
1	B	368	ASN	N-CA-C	5.72	118.29	111.71
1	D	386	VAL	CB-CA-C	-5.33	108.70	114.35
1	E	265	GLY	N-CA-C	5.19	120.59	110.73
1	B	386	VAL	CB-CA-C	-5.17	108.87	114.35
1	E	482	GLY	CA-C-N	5.14	124.93	119.28
1	E	482	GLY	C-N-CA	5.14	124.93	119.28
1	C	386	VAL	CB-CA-C	-5.11	108.93	114.35

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	4864	0	4641	41	0
1	B	4923	0	4703	60	0
1	C	4915	0	4692	35	0
1	D	4880	0	4654	63	0
1	E	4889	0	4667	42	0
1	F	4915	0	4692	47	0
2	A	11	0	0	0	0
2	B	11	0	0	0	0
2	C	11	0	0	0	0
2	D	11	0	0	1	0
2	E	11	0	0	1	0
2	F	11	0	0	1	0
3	A	58	0	0	2	0
3	B	64	0	0	1	0
3	C	72	0	0	1	0
3	D	59	0	0	1	0
3	E	59	0	0	3	0
3	F	66	0	0	1	0
All	All	29830	0	28049	274	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 5.

All (274) 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:248:VAL:HG23	1:B:288:LEU:HD22	1.37	1.03
1:B:248:VAL:HG23	1:B:288:LEU:CD2	1.98	0.92
1:D:82:GLU:HG2	1:F:82:GLU:HG2	1.51	0.89
1:D:263:GLY:HA3	1:D:264:ASN:HB2	1.57	0.87
1:B:16:MET:HE1	1:B:86:LYS:HE3	1.57	0.86
1:B:248:VAL:CG2	1:B:288:LEU:HD22	2.07	0.84
1:F:600:ASN:ND2	1:F:601:GLU:HG2	1.92	0.83
1:B:16:MET:HE2	1:B:182:ALA:HB3	1.63	0.81
1:B:262:ASP:HB2	1:B:264:ASN:H	1.48	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:374:SER:H	1:D:375:GLY:HA2	1.47	0.78
1:B:16:MET:HE1	1:B:86:LYS:CE	2.15	0.76
1:D:277:ARG:O	1:D:277:ARG:HD3	1.84	0.75
1:B:16:MET:HE1	1:B:86:LYS:CD	2.17	0.75
1:B:98:ARG:NH2	1:B:141:GLU:O	2.21	0.74
1:C:155:LEU:HD12	1:C:159:ARG:HG2	1.68	0.74
1:E:257:VAL:HG22	1:E:271:ARG:HG2	1.67	0.74
1:B:16:MET:CE	1:B:182:ALA:HB3	2.17	0.74
1:F:528:MET:O	3:F:801:HOH:O	2.04	0.74
1:D:257:VAL:HG22	1:D:271:ARG:HG2	1.70	0.73
1:D:280:ARG:NH1	1:D:290:ASP:OD1	2.20	0.72
1:F:98:ARG:NH2	1:F:141:GLU:O	2.21	0.72
1:D:567:GLY:HA2	1:E:528:MET:SD	2.30	0.72
1:A:528:MET:O	3:A:801:HOH:O	2.08	0.71
1:A:98:ARG:NH2	1:A:141:GLU:O	2.23	0.71
1:D:156:ASN:ND2	1:E:490:GLU:OE1	2.22	0.71
1:C:16:MET:HE1	1:C:86:LYS:HD2	1.73	0.71
1:E:98:ARG:NH2	1:E:141:GLU:O	2.22	0.70
1:D:277:ARG:HD3	1:D:277:ARG:C	2.17	0.70
1:E:205:LYS:NZ	3:E:802:HOH:O	2.25	0.69
1:C:98:ARG:NH2	1:C:141:GLU:O	2.26	0.68
1:D:98:ARG:NH2	1:D:141:GLU:O	2.25	0.68
1:D:373:GLY:H	1:D:374:SER:HB3	1.59	0.68
1:A:549:GLU:H	1:A:549:GLU:CD	2.04	0.66
1:B:262:ASP:N	1:B:263:GLY:HA2	2.10	0.66
1:D:263:GLY:HA3	1:D:264:ASN:CB	2.27	0.65
1:C:277:ARG:HH11	1:C:279:VAL:HG23	1.62	0.65
1:A:277:ARG:HH11	1:A:279:VAL:HG23	1.62	0.64
1:D:567:GLY:CA	1:E:528:MET:SD	2.85	0.64
1:C:528:MET:O	1:C:529:TRP:HB2	1.98	0.63
1:D:569:MET:HG2	1:E:569:MET:HG2	1.81	0.63
1:A:370:VAL:HG11	1:A:569:MET:HE1	1.81	0.62
1:A:257:VAL:HG22	1:A:271:ARG:HG2	1.81	0.62
1:F:257:VAL:HG22	1:F:271:ARG:HG2	1.81	0.62
1:C:370:VAL:HG21	1:C:569:MET:HE1	1.81	0.61
1:B:549:GLU:CD	1:B:549:GLU:H	2.07	0.61
1:D:338:TYR:O	1:D:403:ARG:NH2	2.29	0.61
1:D:528:MET:O	1:D:529:TRP:HB2	2.00	0.61
1:B:289:ASN:O	1:B:290:ASP:HB2	2.01	0.61
1:B:284:THR:O	1:B:552:GLN:HG3	2.01	0.61
1:B:262:ASP:HB2	1:B:264:ASN:N	2.15	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:GLU:N	1:C:82:GLU:OE1	2.32	0.60
1:D:489:GLU:OE2	1:D:493:ARG:NH1	2.33	0.60
1:D:571:VAL:HG21	1:E:571:VAL:HG21	1.83	0.60
1:F:3:GLU:HA	1:F:189:LYS:NZ	2.17	0.59
1:D:263:GLY:CA	1:D:264:ASN:HB2	2.29	0.59
1:B:82:GLU:OE1	1:B:82:GLU:N	2.35	0.58
1:D:453:LYS:HD2	1:D:479:ILE:HD11	1.83	0.58
1:B:370:VAL:HG21	1:B:569:MET:HE1	1.84	0.58
1:D:372:ALA:HB1	1:D:452:GLU:HG3	1.84	0.58
1:B:248:VAL:O	1:B:249:ARG:HB2	2.04	0.58
1:D:528:MET:SD	1:E:567:GLY:HA2	2.43	0.58
1:A:87:LYS:HE2	1:A:185:GLU:OE1	2.04	0.58
1:C:598:LYS:O	1:C:600:ASN:ND2	2.36	0.58
1:A:289:ASN:O	1:A:290:ASP:HB2	2.03	0.57
1:E:82:GLU:OE1	1:E:82:GLU:N	2.33	0.57
1:B:480:SER:HB2	1:B:488:ALA:HB2	1.86	0.57
1:D:374:SER:H	1:D:375:GLY:CA	2.16	0.57
1:E:289:ASN:O	1:E:290:ASP:HB2	2.03	0.57
1:B:248:VAL:HG22	1:B:352:LEU:HD21	1.87	0.56
1:B:257:VAL:HG22	1:B:271:ARG:HG2	1.87	0.56
1:C:338:TYR:O	1:C:403:ARG:NH2	2.28	0.56
1:A:338:TYR:O	1:A:403:ARG:NH2	2.28	0.56
1:B:452:GLU:HG3	1:B:454:ASN:H	1.71	0.56
1:F:157:ASN:HD21	1:F:159:ARG:HD3	1.71	0.56
1:A:370:VAL:HG11	1:A:569:MET:CE	2.37	0.55
1:A:230:GLU:HG2	1:F:291:ARG:HD2	1.88	0.55
1:E:219:ARG:NH1	3:E:804:HOH:O	2.37	0.55
1:F:480:SER:HB2	1:F:488:ALA:HB2	1.89	0.54
1:D:16:MET:HE1	1:D:86:LYS:HD2	1.90	0.54
1:E:248:VAL:O	1:E:249:ARG:HB2	2.07	0.54
1:F:280:ARG:NH1	1:F:290:ASP:OD1	2.41	0.53
1:C:340:GLU:HG3	1:C:403:ARG:HH21	1.73	0.53
1:D:492:PHE:HD2	1:D:540:MET:HE1	1.74	0.53
1:A:116:LEU:HD13	1:A:399:GLU:HB2	1.90	0.53
1:A:359:ALA:HB3	1:A:416:PHE:HA	1.91	0.53
1:B:571:VAL:O	1:B:575:HIS:NE2	2.42	0.53
1:C:257:VAL:HG22	1:C:271:ARG:HG2	1.91	0.52
1:E:280:ARG:NH1	1:E:290:ASP:OD1	2.39	0.52
1:D:502:LYS:HB3	1:D:504:LEU:HD13	1.92	0.52
1:B:16:MET:CE	1:B:86:LYS:CD	2.86	0.52
1:D:596:TRP:O	1:D:597:GLU:HB2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:MET:HB2	1:B:182:ALA:HB3	1.92	0.52
1:B:5:SER:OG	1:B:189:LYS:O	2.22	0.52
1:B:6:GLU:HG2	1:B:403:ARG:HB2	1.91	0.51
1:D:528:MET:O	1:D:529:TRP:CB	2.58	0.51
1:B:63:LYS:HG3	1:B:159:ARG:HH12	1.75	0.51
1:C:599:LYS:NZ	3:C:802:HOH:O	2.39	0.51
1:B:249:ARG:O	1:B:250:ASN:HB2	2.11	0.51
1:D:116:LEU:HD13	1:D:399:GLU:HB2	1.92	0.51
1:D:370:VAL:HB	1:D:569:MET:HE1	1.93	0.51
1:C:297:GLY:HA3	1:C:330:CYS:O	2.11	0.51
1:E:82:GLU:CD	1:E:82:GLU:H	2.14	0.51
1:A:207:GLU:HG2	1:A:235:GLU:HG2	1.94	0.50
1:D:262:ASP:HB3	1:D:267:LEU:HD11	1.92	0.50
1:C:203:LEU:HD22	1:C:239:GLN:HG2	1.94	0.50
1:D:568:ILE:HG13	1:D:569:MET:HG3	1.93	0.50
1:F:296:LYS:NZ	1:F:599:LYS:O	2.38	0.50
1:D:418:GLU:HG2	1:D:452:GLU:HB2	1.93	0.49
1:C:204:VAL:HB	1:C:238:LEU:HB2	1.94	0.49
1:E:363:MET:HE3	1:E:365:SER:HB3	1.94	0.49
1:A:203:LEU:HD21	1:F:244:ARG:HD3	1.94	0.49
1:B:458:GLU:OE2	1:B:498:ARG:HD2	2.12	0.49
1:D:289:ASN:O	1:D:290:ASP:HB2	2.11	0.49
1:B:568:ILE:HG13	1:B:569:MET:HG3	1.95	0.49
1:A:6:GLU:HG2	1:A:403:ARG:HB2	1.95	0.49
1:A:480:SER:HB2	1:A:488:ALA:HB2	1.93	0.49
1:D:334:SER:HA	1:D:335:HIS:HA	1.68	0.48
1:B:82:GLU:CD	1:B:82:GLU:H	2.15	0.48
1:E:186:GLU:OE1	1:E:214:HIS:ND1	2.37	0.48
1:A:145:THR:HG22	1:A:385:THR:HB	1.95	0.48
1:B:246:TRP:CZ2	1:B:350:GLY:HA2	2.49	0.48
1:D:374:SER:N	1:D:375:GLY:HA2	2.17	0.48
1:E:282:GLU:OE1	3:E:801:HOH:O	2.20	0.48
1:C:246:TRP:CZ2	1:C:350:GLY:HA2	2.48	0.48
1:B:16:MET:CE	1:B:182:ALA:CB	2.91	0.47
1:B:247:GLU:HB3	1:B:250:ASN:HB3	1.96	0.47
1:E:297:GLY:HA3	1:E:330:CYS:O	2.15	0.47
1:E:475:TYR:O	1:E:540:MET:HE1	2.15	0.47
1:C:96:THR:HA	1:C:97:HIS:HA	1.69	0.47
1:C:277:ARG:HG3	1:C:288:LEU:HD11	1.96	0.47
1:C:489:GLU:O	1:C:493:ARG:HG3	2.15	0.47
1:D:204:VAL:HB	1:D:238:LEU:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:332:ARG:NH1	1:D:356:GLU:OE2	2.46	0.47
1:F:418:GLU:HG2	1:F:452:GLU:HB2	1.97	0.47
1:A:297:GLY:HA3	1:A:330:CYS:O	2.14	0.47
1:F:477:TRP:CZ2	1:F:516:ASP:HB2	2.50	0.47
1:F:141:GLU:OE1	1:F:160:LYS:NZ	2.45	0.47
1:F:600:ASN:CG	1:F:601:GLU:N	2.73	0.47
1:F:489:GLU:O	1:F:493:ARG:HG3	2.14	0.47
1:C:458:GLU:OE2	1:C:498:ARG:HD2	2.14	0.46
1:D:448:THR:OG1	1:D:449:GLY:N	2.47	0.46
1:F:248:VAL:O	1:F:249:ARG:HB2	2.16	0.46
1:A:96:THR:HA	1:A:97:HIS:HA	1.71	0.46
1:A:219:ARG:NH2	1:F:600:ASN:HD21	2.14	0.46
1:C:557:TRP:HA	1:C:558:ASN:HA	1.70	0.46
1:E:489:GLU:O	1:E:493:ARG:HG3	2.14	0.46
1:E:448:THR:OG1	1:E:449:GLY:N	2.48	0.46
1:E:369:PHE:N	1:E:369:PHE:CD1	2.81	0.46
1:B:448:THR:OG1	1:B:449:GLY:N	2.48	0.46
1:D:493:ARG:NH2	3:D:804:HOH:O	2.48	0.46
1:E:458:GLU:OE2	1:E:498:ARG:HD2	2.15	0.46
1:E:334:SER:HA	1:E:335:HIS:HA	1.60	0.46
1:A:448:THR:OG1	1:A:449:GLY:N	2.48	0.46
1:D:435:ALA:O	1:D:439:THR:HG23	2.16	0.46
1:A:277:ARG:NH1	1:A:279:VAL:HG23	2.31	0.45
1:B:16:MET:HE2	1:B:16:MET:HB2	1.70	0.45
1:E:116:LEU:HD13	1:E:399:GLU:HB2	1.98	0.45
1:B:578:VAL:HG12	1:B:579:PHE:CD2	2.50	0.45
1:E:96:THR:HA	1:E:97:HIS:HA	1.71	0.45
1:F:206:TYR:CE1	1:F:218:VAL:HG21	2.50	0.45
1:A:458:GLU:OE2	1:A:498:ARG:HD2	2.17	0.45
1:F:297:GLY:HA3	1:F:330:CYS:O	2.17	0.45
1:F:600:ASN:C	1:F:601:GLU:CG	2.88	0.45
1:A:369:PHE:CD1	1:A:370:VAL:HG13	2.52	0.45
1:B:477:TRP:CZ2	1:B:516:ASP:HB2	2.52	0.45
1:C:119:LEU:HA	1:C:119:LEU:HD23	1.77	0.45
1:C:453:LYS:HD2	1:C:479:ILE:HD11	1.97	0.45
1:F:571:VAL:O	1:F:575:HIS:NE2	2.49	0.45
1:A:378:THR:OG1	1:A:420:GLU:OE2	2.32	0.45
1:C:277:ARG:HE	1:C:408:PRO:HA	1.82	0.45
1:C:418:GLU:HG2	1:C:452:GLU:HB2	1.99	0.45
1:D:312:HIS:CD2	1:F:312:HIS:HD2	2.35	0.45
1:D:23:MET:HE3	1:F:345:PHE:CE1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:386:VAL:HG22	1:A:387:PRO:HD3	1.97	0.45
1:B:470:CYS:HA	1:B:509:VAL:O	2.18	0.44
1:C:352:LEU:HB3	1:C:411:ILE:HD13	1.97	0.44
1:C:334:SER:HA	1:C:335:HIS:HA	1.67	0.44
1:D:453:LYS:HD2	1:D:479:ILE:CD1	2.47	0.44
1:E:204:VAL:HB	1:E:238:LEU:HB2	2.00	0.44
1:E:167:ASP:OD2	2:E:701:SJ5:O4	2.35	0.44
1:B:96:THR:HA	1:B:97:HIS:HA	1.66	0.44
1:F:155:LEU:HD23	1:F:159:ARG:HG2	2.00	0.44
1:F:457:PRO:O	1:F:502:LYS:NZ	2.47	0.44
1:A:470:CYS:HA	1:A:509:VAL:O	2.17	0.44
1:A:204:VAL:HB	1:A:238:LEU:HB2	2.00	0.44
1:A:284:THR:O	1:A:552:GLN:HG3	2.18	0.44
1:D:528:MET:O	1:D:529:TRP:CD1	2.71	0.44
1:F:334:SER:HA	1:F:335:HIS:HA	1.63	0.44
1:B:16:MET:HE1	1:B:86:LYS:HD3	1.98	0.44
1:C:116:LEU:HD13	1:C:399:GLU:HB2	1.99	0.44
1:D:207:GLU:HG2	1:D:235:GLU:HG2	2.00	0.44
1:B:116:LEU:HD13	1:B:399:GLU:HB2	2.00	0.43
1:B:334:SER:HA	1:B:335:HIS:HA	1.67	0.43
1:B:489:GLU:O	1:B:493:ARG:HG3	2.18	0.43
1:B:206:TYR:CE1	1:B:218:VAL:HG21	2.53	0.43
1:B:477:TRP:HB2	1:B:515:THR:HG22	2.00	0.43
1:B:557:TRP:HA	1:B:558:ASN:HA	1.72	0.43
1:C:435:ALA:O	1:C:439:THR:HG23	2.18	0.43
1:A:334:SER:HA	1:A:335:HIS:HA	1.60	0.43
1:A:452:GLU:HG3	1:A:454:ASN:H	1.82	0.43
1:D:3:GLU:O	1:D:3:GLU:HG2	2.18	0.43
1:A:557:TRP:HA	1:A:558:ASN:HA	1.72	0.43
1:C:359:ALA:HB3	1:C:416:PHE:HA	2.00	0.43
1:D:297:GLY:HA3	1:D:330:CYS:O	2.18	0.43
1:A:247:GLU:HB2	1:A:250:ASN:HB3	2.00	0.43
1:F:568:ILE:HG13	1:F:569:MET:HG3	2.00	0.43
1:B:297:GLY:HA3	1:B:330:CYS:O	2.19	0.43
1:C:82:GLU:H	1:C:82:GLU:CD	2.26	0.43
1:E:277:ARG:HG3	1:E:288:LEU:HD11	2.01	0.43
1:F:82:GLU:H	1:F:82:GLU:CD	2.27	0.43
1:B:190:ASP:OD1	3:B:801:HOH:O	2.22	0.42
1:D:359:ALA:HB3	1:D:416:PHE:HA	2.00	0.42
1:F:362:MET:HE1	1:F:393:HIS:ND1	2.33	0.42
1:B:145:THR:HG22	1:B:385:THR:HB	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:463:TYR:CG	1:B:464:PRO:HD3	2.54	0.42
1:C:206:TYR:CE1	1:C:218:VAL:HG21	2.53	0.42
1:F:222:ASP:HB2	1:F:252:TYR:OH	2.19	0.42
1:F:359:ALA:HB3	1:F:416:PHE:HA	2.02	0.42
1:B:578:VAL:HG21	1:B:592:PHE:CE2	2.53	0.42
1:A:332:ARG:NH1	1:A:356:GLU:OE2	2.52	0.42
1:A:489:GLU:O	1:A:493:ARG:HG3	2.18	0.42
1:D:373:GLY:N	1:D:374:SER:HB3	2.32	0.42
1:B:587:ALA:O	1:B:590:VAL:HG22	2.19	0.42
1:E:164:PRO:HB3	1:E:166:PHE:CZ	2.55	0.42
1:E:284:THR:O	1:E:552:GLN:HG3	2.20	0.42
1:B:363:MET:HE3	1:B:365:SER:HB3	2.02	0.42
1:C:470:CYS:HA	1:C:509:VAL:O	2.19	0.42
1:D:36:LYS:HB3	1:D:36:LYS:HE2	1.96	0.42
1:F:486:GLU:O	1:F:490:GLU:HG2	2.19	0.42
1:C:334:SER:HB3	1:C:335:HIS:CE1	2.54	0.42
1:D:335:HIS:O	1:D:358:PRO:HA	2.20	0.42
1:F:557:TRP:HA	1:F:558:ASN:HA	1.71	0.42
1:D:277:ARG:C	1:D:277:ARG:CD	2.86	0.42
1:E:244:ARG:HB3	1:E:252:TYR:CD2	2.55	0.42
1:A:335:HIS:O	1:A:358:PRO:HA	2.20	0.42
1:E:557:TRP:HA	1:E:558:ASN:HA	1.72	0.42
1:E:470:CYS:HA	1:E:509:VAL:O	2.20	0.41
1:B:16:MET:HE3	1:B:182:ALA:HB3	1.98	0.41
1:D:512:GLU:OE1	2:D:701:SJ5:N	2.53	0.41
1:D:206:TYR:CE1	1:D:218:VAL:HG21	2.55	0.41
1:F:23:MET:HE3	1:F:48:SER:HB3	2.02	0.41
1:B:335:HIS:O	1:B:358:PRO:HA	2.20	0.41
1:B:475:TYR:O	1:B:540:MET:HE1	2.20	0.41
1:E:304:PHE:CG	1:E:305:PRO:HD2	2.56	0.41
1:A:463:TYR:CG	1:A:464:PRO:HD3	2.56	0.41
1:D:82:GLU:HG2	1:F:82:GLU:CG	2.38	0.41
1:F:111:HIS:CG	1:F:118:VAL:HG22	2.55	0.41
1:F:289:ASN:O	1:F:290:ASP:HB2	2.20	0.41
1:C:475:TYR:O	1:C:540:MET:HE1	2.19	0.41
1:A:12:ASN:HB2	3:A:843:HOH:O	2.21	0.41
1:D:18:GLN:CD	1:F:11:GLN:HE22	2.28	0.41
1:A:475:TYR:O	1:A:540:MET:HE1	2.21	0.41
1:D:528:MET:O	1:D:529:TRP:HD1	2.04	0.41
1:E:367:ARG:O	1:E:367:ARG:HG3	2.21	0.41
1:F:597:GLU:HG2	1:F:598:LYS:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:ILE:HG21	1:B:120:ALA:HB1	2.03	0.41
1:D:418:GLU:CG	1:D:452:GLU:HB2	2.51	0.41
1:E:108:ILE:HG21	1:E:120:ALA:HB1	2.03	0.41
1:F:14:TYR:O	1:F:184:PRO:HD3	2.21	0.41
1:A:280:ARG:NH1	1:A:282:GLU:OE2	2.53	0.40
1:E:246:TRP:CZ2	1:E:350:GLY:HA2	2.56	0.40
1:E:255:GLN:OE1	1:E:271:ARG:HD3	2.20	0.40
1:F:304:PHE:CG	1:F:305:PRO:HD2	2.56	0.40
1:E:206:TYR:CE1	1:E:218:VAL:HG21	2.56	0.40
1:D:463:TYR:CG	1:D:464:PRO:HD3	2.56	0.40
1:D:470:CYS:HA	1:D:509:VAL:O	2.21	0.40
1:E:277:ARG:NH2	1:E:443:GLN:OE1	2.51	0.40
1:F:470:CYS:HA	1:F:509:VAL:O	2.21	0.40
1:F:512:GLU:OE1	2:F:701:SJ5:N	2.54	0.40
1:B:277:ARG:HG3	1:B:288:LEU:HD11	2.04	0.40
1:D:111:HIS:CG	1:D:118:VAL:HG22	2.57	0.40
1:F:452:GLU:HG2	1:F:453:LYS:H	1.86	0.40
1:D:571:VAL:O	1:D:575:HIS:NE2	2.55	0.40
1:F:458:GLU:OE2	1:F:498:ARG:HD2	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	589/627 (94%)	571 (97%)	17 (3%)	1 (0%)	43	63
1	B	599/627 (96%)	577 (96%)	20 (3%)	2 (0%)	36	55
1	C	598/627 (95%)	578 (97%)	20 (3%)	0	100	100
1	D	594/627 (95%)	570 (96%)	21 (4%)	3 (0%)	24	43
1	E	595/627 (95%)	575 (97%)	19 (3%)	1 (0%)	43	63

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	598/627 (95%)	578 (97%)	20 (3%)	0	100	100
All	All	3573/3762 (95%)	3449 (96%)	117 (3%)	7 (0%)	43	63

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	374	SER
1	D	264	ASN
1	D	529	TRP
1	B	249	ARG
1	B	250	ASN
1	D	376	ASN
1	E	249	ARG

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	517/546 (95%)	513 (99%)	4 (1%)	73	88
1	B	523/546 (96%)	511 (98%)	12 (2%)	44	72
1	C	522/546 (96%)	517 (99%)	5 (1%)	68	86
1	D	518/546 (95%)	507 (98%)	11 (2%)	47	74
1	E	519/546 (95%)	511 (98%)	8 (2%)	57	80
1	F	522/546 (96%)	510 (98%)	12 (2%)	44	72
All	All	3121/3276 (95%)	3069 (98%)	52 (2%)	53	78

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	LYS
1	A	134	VAL
1	A	244	ARG
1	A	277	ARG

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Mol	Chain	Res	Type
1	B	6	GLU
1	B	82	GLU
1	B	119	LEU
1	B	134	VAL
1	B	156	ASN
1	B	277	ARG
1	B	334	SER
1	B	367	ARG
1	B	370	VAL
1	B	391	LYS
1	B	549	GLU
1	B	578	VAL
1	C	35	LYS
1	C	159	ARG
1	C	280	ARG
1	C	334	SER
1	C	370	VAL
1	D	36	LYS
1	D	63	LYS
1	D	134	VAL
1	D	224	GLU
1	D	232	GLU
1	D	244	ARG
1	D	277	ARG
1	D	334	SER
1	D	521	LEU
1	D	540	MET
1	D	597	GLU
1	E	82	GLU
1	E	134	VAL
1	E	277	ARG
1	E	369	PHE
1	E	370	VAL
1	E	384	LEU
1	E	424	ASP
1	E	528	MET
1	F	119	LEU
1	F	155	LEU
1	F	244	ARG
1	F	248	VAL
1	F	264	ASN
1	F	277	ARG

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Mol	Chain	Res	Type
1	F	279	VAL
1	F	490	GLU
1	F	578	VAL
1	F	597	GLU
1	F	600	ASN
1	F	601	GLU

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

Mol	Chain	Res	Type
1	A	312	HIS
1	B	11	GLN
1	B	443	GLN
1	C	552	GLN
1	C	584	GLN
1	D	11	GLN
1	E	11	GLN
1	E	130	GLN
1	E	327	ASN
1	E	552	GLN
1	F	11	GLN
1	F	563	GLN
1	F	600	ASN

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](#)

6 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	$\# Z > 2$
2	SJ5	E	701	-	11,11,11	0.51	0	11,15,15	2.99	3 (27%)
2	SJ5	D	701	-	11,11,11	0.51	0	11,15,15	2.80	3 (27%)
2	SJ5	C	701	-	11,11,11	0.49	0	11,15,15	2.80	3 (27%)
2	SJ5	A	701	-	11,11,11	0.50	0	11,15,15	3.27	3 (27%)
2	SJ5	F	701	-	11,11,11	0.45	0	11,15,15	2.66	3 (27%)
2	SJ5	B	701	-	11,11,11	0.49	0	11,15,15	2.78	3 (27%)

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	SJ5	E	701	-	-	2/4/18/18	0/1/1/1
2	SJ5	D	701	-	-	2/4/18/18	0/1/1/1
2	SJ5	C	701	-	-	2/4/18/18	0/1/1/1
2	SJ5	A	701	-	-	2/4/18/18	0/1/1/1
2	SJ5	F	701	-	-	2/4/18/18	0/1/1/1
2	SJ5	B	701	-	-	2/4/18/18	0/1/1/1

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	SJ5	C7-N-C2	10.29	123.71	111.78
2	E	701	SJ5	C7-N-C2	9.29	122.55	111.78
2	C	701	SJ5	C7-N-C2	8.61	121.76	111.78
2	D	701	SJ5	C7-N-C2	8.52	121.65	111.78
2	B	701	SJ5	C7-N-C2	8.51	121.65	111.78
2	F	701	SJ5	C7-N-C2	8.02	121.08	111.78
2	D	701	SJ5	O62-C6-C5	2.54	120.52	113.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	701	SJ5	O62-C6-C5	2.49	120.40	113.93
2	B	701	SJ5	O62-C6-C5	2.49	120.40	113.93
2	A	701	SJ5	O62-C6-C5	2.41	120.17	113.93
2	C	701	SJ5	O62-C6-C5	2.37	120.07	113.93
2	E	701	SJ5	O62-C6-C5	2.36	120.05	113.93
2	B	701	SJ5	O62-C6-O61	-2.23	119.01	124.08
2	D	701	SJ5	O62-C6-O61	-2.23	119.02	124.08
2	E	701	SJ5	O62-C6-O61	-2.21	119.06	124.08
2	C	701	SJ5	O62-C6-O61	-2.19	119.11	124.08
2	A	701	SJ5	O62-C6-O61	-2.16	119.17	124.08
2	F	701	SJ5	O62-C6-O61	-2.14	119.23	124.08

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	701	SJ5	C4-C5-C6-O62
2	B	701	SJ5	C4-C5-C6-O61
2	E	701	SJ5	C4-C5-C6-O62
2	C	701	SJ5	C4-C5-C6-O61
2	D	701	SJ5	C4-C5-C6-O61
2	E	701	SJ5	C4-C5-C6-O61
2	F	701	SJ5	C4-C5-C6-O62
2	F	701	SJ5	C4-C5-C6-O61
2	A	701	SJ5	C4-C5-C6-O61
2	A	701	SJ5	C4-C5-C6-O62
2	C	701	SJ5	C4-C5-C6-O62
2	D	701	SJ5	C4-C5-C6-O62

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	701	SJ5	1	0
2	D	701	SJ5	1	0
2	F	701	SJ5	1	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	593/627 (94%)	-0.16	8 (1%) 75 71	27, 39, 63, 100	0
1	B	601/627 (95%)	-0.17	6 (0%) 79 76	27, 40, 62, 95	0
1	C	600/627 (95%)	-0.20	1 (0%) 91 89	27, 41, 64, 92	0
1	D	596/627 (95%)	-0.27	5 (0%) 82 80	26, 38, 63, 100	0
1	E	597/627 (95%)	-0.22	7 (1%) 76 73	28, 39, 65, 95	0
1	F	600/627 (95%)	-0.21	5 (0%) 82 80	27, 40, 62, 93	0
All	All	3587/3762 (95%)	-0.21	32 (0%) 81 78	26, 39, 64, 100	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	249	ARG	5.3
1	E	528	MET	4.4
1	E	369	PHE	3.8
1	F	528	MET	3.6
1	F	2	LEU	3.4
1	E	290	ASP	3.2
1	E	263	GLY	3.2
1	A	373	GLY	3.1
1	A	481	GLY	3.1
1	B	248	VAL	2.7
1	A	290	ASP	2.6
1	A	369	PHE	2.6
1	A	549	GLU	2.5
1	B	549	GLU	2.4
1	B	290	ASP	2.4
1	D	528	MET	2.4
1	B	372	ALA	2.4
1	C	156	ASN	2.3
1	F	290	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	549	GLU	2.3
1	E	264	ASN	2.2
1	A	262	ASP	2.2
1	F	3	GLU	2.2
1	E	373	GLY	2.1
1	F	264	ASN	2.1
1	A	375	GLY	2.1
1	D	263	GLY	2.1
1	D	372	ALA	2.0
1	A	452	GLU	2.0
1	B	85	ASN	2.0
1	D	264	ASN	2.0
1	D	375	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
2	SJ5	B	701	11/11	0.92	0.08	34,38,50,54	0
2	SJ5	C	701	11/11	0.93	0.07	31,35,38,40	0
2	SJ5	E	701	11/11	0.94	0.08	32,34,38,43	0
2	SJ5	F	701	11/11	0.94	0.09	35,39,42,43	0
2	SJ5	A	701	11/11	0.96	0.07	31,37,40,43	0
2	SJ5	D	701	11/11	0.96	0.08	35,37,42,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.