



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:27 AM UTC

PDB ID : 3QTD / pdb_00003qtd
Title : Crystal structure of putative modulator of gyrase (PmbA) from *Pseudomonas aeruginosa* PAO1
Authors : Tkaczuk, K.L.; Chruszcz, M.; Evdokimova, E.; Liu, F.; Savchenko, A.; Edwards, A.; Joachimiak, A.; Minor, W.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2011-02-22
Resolution : 2.70 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

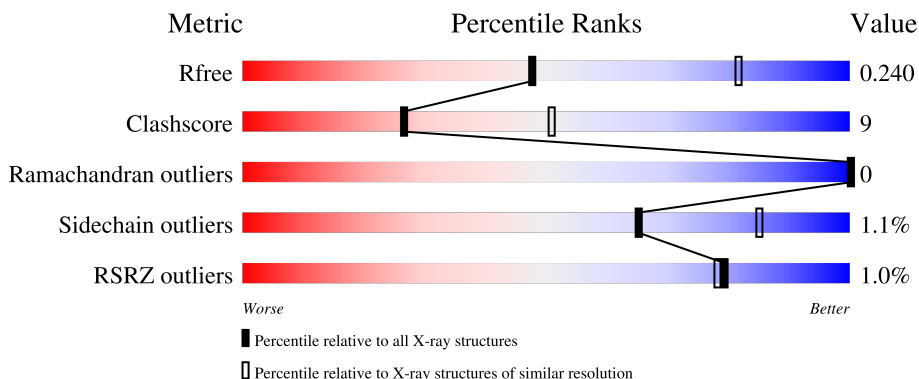
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	449	71% 24% ..
1	B	449	75% 22% ..
1	C	449	2% 81% 16% ..
1	D	449	77% 19% ..

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	GOL	D	451	-	-	X	-

2 Entry composition [i](#)

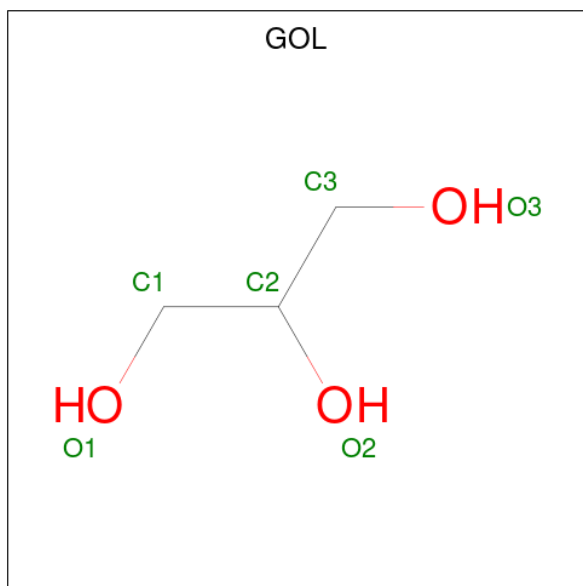
There are 3 unique types of molecules in this entry. The entry contains 13604 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PmbA protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	Se	0	0	0
			3298	2048	597	641	5	7			
1	B	441	Total	C	N	O	S	Se	0	0	0
			3303	2050	600	641	5	7			
1	C	442	Total	C	N	O	S	Se	0	0	0
			3295	2047	599	637	5	7			
1	D	441	Total	C	N	O	S	Se	0	0	0
			3299	2048	600	639	5	7			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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

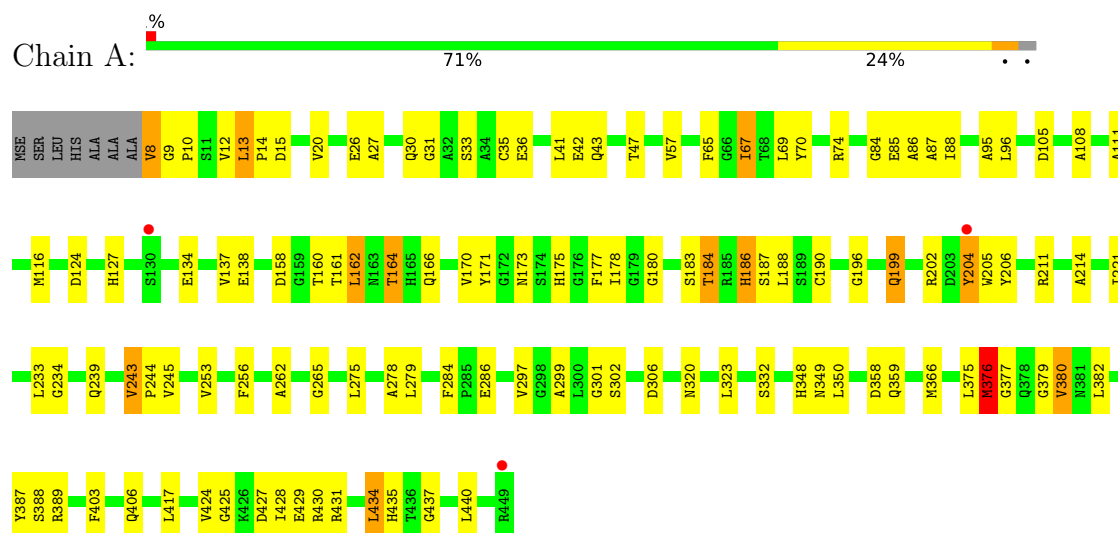
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	86	Total O 86 86	0	0
3	B	112	Total O 112 112	0	0
3	C	98	Total O 98 98	0	0
3	D	59	Total O 59 59	0	0

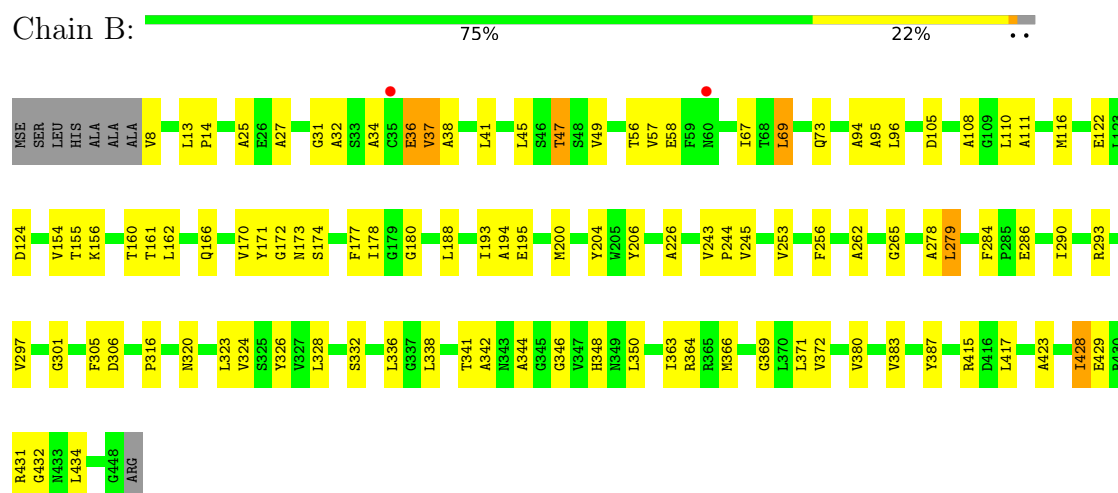
3 Residue-property plots

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.

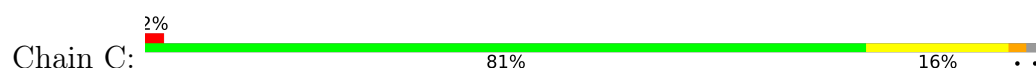
• Molecule 1: PmbA protein

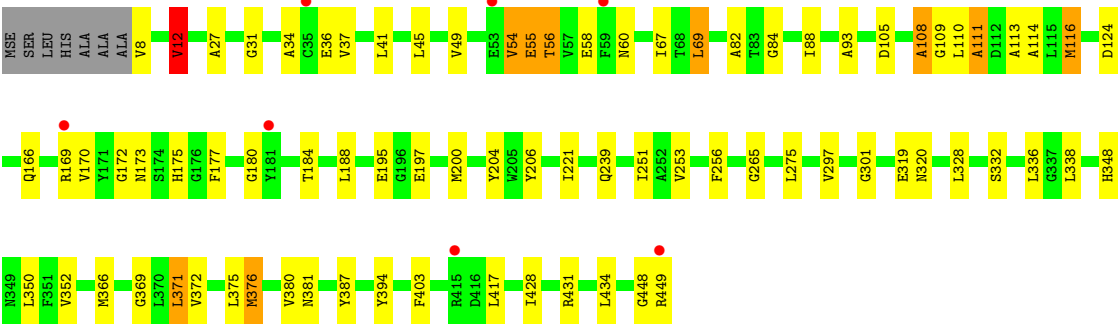


• Molecule 1: PmbA protein

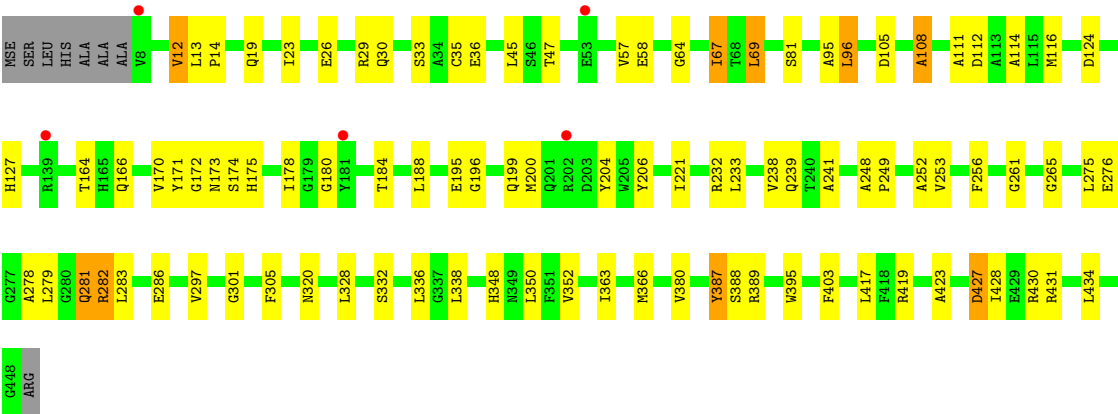
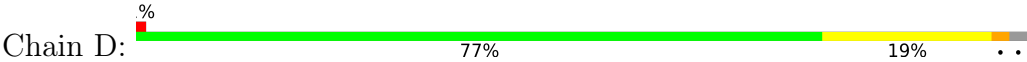


• Molecule 1: PmbA protein





● Molecule 1: PmbA protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.39Å 116.20Å 253.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 50.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.5 (50.00-2.70) 97.5 (50.00-2.70)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	0.17	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.29 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.6.0070	Depositor
R, R_{free}	0.225 , 0.279 (Not available) , 0.240	Depositor DCC
R_{free} test set	2831 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	33.3	Xtriage
Anisotropy	0.808	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13604	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.74 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.1420e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

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

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	1.58	32/3347 (1.0%)	1.14	10/4522 (0.2%)
1	B	1.60	32/3352 (1.0%)	1.14	5/4527 (0.1%)
1	C	1.34	20/3344 (0.6%)	1.09	9/4517 (0.2%)
1	D	1.34	16/3348 (0.5%)	1.09	7/4522 (0.2%)
All	All	1.47	100/13391 (0.7%)	1.12	31/18088 (0.2%)

All (100) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	342	ALA	CA-CB	-8.26	1.41	1.53
1	B	290	ILE	C-O	-7.67	1.16	1.24
1	B	284	PHE	C-O	-7.63	1.17	1.24
1	B	316	PRO	CA-C	7.41	1.60	1.52
1	B	305	PHE	CA-C	-7.39	1.43	1.52
1	B	344	ALA	CA-CB	-7.34	1.43	1.53
1	B	432	GLY	C-O	-7.32	1.18	1.24
1	C	82	ALA	CA-CB	-7.28	1.41	1.53
1	B	49	VAL	C-O	-7.18	1.16	1.24
1	A	245	VAL	C-O	-6.85	1.17	1.24
1	D	67	ILE	C-O	-6.84	1.17	1.24
1	C	111	ALA	CA-CB	-6.81	1.42	1.53
1	B	34	ALA	CA-CB	-6.70	1.42	1.53
1	A	204	TYR	C-O	-6.54	1.16	1.23
1	A	67	ILE	C-O	-6.50	1.17	1.24
1	A	162	LEU	C-O	-6.39	1.16	1.24
1	A	86	ALA	CA-CB	-6.34	1.43	1.53
1	B	47	THR	C-O	-6.32	1.16	1.24
1	B	245	VAL	C-O	-6.26	1.17	1.24
1	A	85	GLU	C-O	-6.25	1.16	1.24
1	B	306	ASP	C-O	-6.23	1.15	1.23
1	B	37	VAL	C-O	-6.19	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	306	ASP	C-O	-6.16	1.16	1.23
1	A	183	SER	C-O	-6.13	1.16	1.23
1	C	27	ALA	CA-CB	-6.11	1.43	1.53
1	A	161	THR	C-O	-6.02	1.16	1.24
1	A	187	SER	C-O	-5.93	1.16	1.23
1	C	116	MSE	SE-CE	-5.92	1.77	1.95
1	D	283	LEU	C-O	-5.91	1.16	1.24
1	D	305	PHE	C-O	-5.91	1.16	1.23
1	C	169	ARG	C-O	-5.82	1.16	1.23
1	C	114	ALA	CA-CB	-5.81	1.43	1.53
1	B	73	GLN	C-O	-5.78	1.16	1.24
1	A	205	TRP	C-O	-5.76	1.16	1.23
1	A	160	THR	C-O	-5.74	1.17	1.24
1	D	178	ILE	C-O	-5.66	1.18	1.24
1	D	252	ALA	CA-CB	5.65	1.62	1.53
1	B	160	THR	C-O	-5.63	1.17	1.24
1	A	127	HIS	N-CA	-5.63	1.40	1.46
1	B	428	ILE	C-O	-5.61	1.17	1.24
1	C	54	VAL	CA-CB	-5.59	1.47	1.53
1	A	27	ALA	CA-CB	-5.59	1.44	1.53
1	B	56	THR	C-O	-5.58	1.17	1.23
1	B	34	ALA	C-O	-5.58	1.17	1.23
1	A	214	ALA	CA-CB	-5.56	1.44	1.53
1	C	37	VAL	C-O	-5.56	1.18	1.24
1	B	194	ALA	C-O	-5.56	1.16	1.23
1	C	116	MSE	CG-SE	-5.55	1.78	1.95
1	A	186	HIS	C-O	-5.52	1.17	1.23
1	C	56	THR	C-O	-5.50	1.17	1.23
1	A	15	ASP	C-O	-5.50	1.17	1.24
1	B	262	ALA	CA-CB	-5.50	1.44	1.53
1	A	349	ASN	CA-C	-5.46	1.48	1.53
1	D	428	ILE	CA-CB	-5.45	1.48	1.54
1	D	172	GLY	C-O	-5.44	1.17	1.24
1	C	34	ALA	C-O	-5.40	1.17	1.23
1	D	427	ASP	C-O	-5.39	1.17	1.23
1	D	114	ALA	CA-CB	-5.39	1.44	1.53
1	A	424	VAL	C-O	-5.38	1.18	1.24
1	A	164	THR	C-O	-5.37	1.17	1.24
1	A	406	GLN	CA-C	-5.36	1.46	1.52
1	A	284	PHE	C-O	-5.35	1.17	1.24
1	D	69	LEU	C-O	-5.32	1.17	1.23
1	C	239	GLN	C-O	-5.29	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	25	ALA	C-O	-5.29	1.18	1.24
1	B	344	ALA	C-O	-5.25	1.17	1.24
1	A	211	ARG	C-O	-5.23	1.17	1.24
1	A	302	SER	C-O	-5.22	1.17	1.23
1	B	293	ARG	C-O	-5.21	1.18	1.24
1	C	428	ILE	CA-CB	-5.21	1.48	1.54
1	B	32	ALA	CA-CB	-5.20	1.44	1.53
1	B	27	ALA	CA-CB	-5.19	1.45	1.53
1	D	241	ALA	CA-CB	-5.19	1.45	1.53
1	A	87	ALA	CA-CB	-5.18	1.45	1.53
1	B	279	LEU	N-CA	-5.18	1.39	1.46
1	B	110	LEU	C-O	-5.17	1.16	1.23
1	B	172	GLY	C-O	-5.15	1.18	1.24
1	A	13	LEU	C-O	-5.14	1.19	1.24
1	B	316	PRO	C-O	5.14	1.29	1.23
1	A	430	ARG	C-O	-5.12	1.17	1.24
1	B	324	VAL	C-O	-5.12	1.16	1.24
1	C	34	ALA	CA-CB	-5.11	1.45	1.53
1	D	281	GLN	CA-C	-5.11	1.46	1.52
1	C	54	VAL	C-O	-5.11	1.17	1.23
1	D	282	ARG	CA-C	-5.08	1.46	1.52
1	C	12	VAL	C-O	-5.08	1.18	1.24
1	A	262	ALA	CA-CB	-5.07	1.44	1.53
1	C	113	ALA	CA-CB	-5.07	1.45	1.53
1	D	112	ASP	C-O	-5.06	1.17	1.23
1	A	184	THR	C-O	-5.04	1.17	1.23
1	B	174	SER	C-O	-5.04	1.17	1.24
1	A	427	ASP	C-O	-5.04	1.18	1.23
1	A	376	MSE	CA-C	-5.03	1.46	1.52
1	C	93	ALA	N-CA	5.03	1.52	1.46
1	B	36	GLU	C-O	-5.03	1.18	1.24
1	C	108	ALA	CA-CB	-5.02	1.45	1.53
1	C	49	VAL	C-O	-5.01	1.18	1.23
1	D	108	ALA	CA-CB	-5.01	1.45	1.53
1	D	387	TYR	C-O	-5.01	1.18	1.23
1	A	380	VAL	C-O	-5.00	1.18	1.24

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	428	ILE	N-CA-C	9.66	121.64	108.89
1	B	428	ILE	N-CA-C	8.26	120.23	109.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	55	GLU	N-CA-C	6.97	118.95	111.36
1	D	403	PHE	N-CA-C	6.95	114.47	108.22
1	C	54	VAL	CB-CA-C	-6.63	102.83	111.25
1	D	428	ILE	N-CA-C	6.59	118.08	109.58
1	D	276	GLU	N-CA-C	6.18	118.01	111.28
1	C	403	PHE	N-CA-C	6.00	113.62	108.22
1	C	428	ILE	N-CA-C	5.86	117.14	109.58
1	A	403	PHE	N-CA-C	5.84	113.38	108.13
1	D	423	ALA	N-CA-C	5.79	117.64	108.79
1	B	173	ASN	N-CA-C	5.78	116.71	108.74
1	A	96	LEU	N-CA-C	5.68	117.55	111.36
1	C	376	MSE	N-CA-C	5.56	117.72	108.99
1	D	430	ARG	N-CA-C	5.54	120.73	113.30
1	A	199	GLN	N-CA-C	5.48	117.13	108.96
1	A	434	LEU	CD1-CG-CD2	-5.46	98.79	110.80
1	B	369	GLY	N-CA-C	5.32	117.19	110.38
1	D	261	GLY	N-CA-C	-5.30	106.37	112.73
1	B	31	GLY	N-CA-C	5.18	122.57	115.27
1	A	31	GLY	N-CA-C	5.15	122.69	115.43
1	A	243	VAL	CA-C-N	5.14	125.04	119.85
1	A	243	VAL	C-N-CA	5.14	125.04	119.85
1	C	109	GLY	N-CA-C	5.13	116.83	111.95
1	C	31	GLY	N-CA-C	5.09	122.61	115.43
1	A	33	SER	CB-CA-C	-5.08	102.22	110.85
1	A	211	ARG	CB-CA-C	-5.06	101.99	110.19
1	B	346	GLY	N-CA-C	5.05	121.91	112.77
1	D	96	LEU	N-CA-C	5.04	116.86	111.36
1	C	369	GLY	N-CA-C	5.04	116.83	110.38
1	C	37	VAL	N-CA-C	5.00	115.91	108.46

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	3298	0	3210	72	0
1	B	3303	0	3225	67	0
1	C	3295	0	3211	49	0
1	D	3299	0	3221	65	0
2	A	12	0	15	0	0
2	B	12	0	16	3	0
2	C	12	0	16	1	0
2	D	18	0	24	5	0
3	A	86	0	0	1	0
3	B	112	0	0	2	0
3	C	98	0	0	2	0
3	D	59	0	0	2	0
All	All	13604	0	12938	243	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (243) 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:C:375:LEU:C	1:C:376:MSE:HG3	1.59	1.15
1:A:8:VAL:HG13	1:A:41:LEU:HD23	1.39	1.00
1:C:375:LEU:O	1:C:376:MSE:HG3	1.62	0.98
1:D:278:ALA:O	1:D:281:GLN:HG3	1.68	0.93
1:B:350:LEU:HD23	1:B:434:LEU:CD1	2.06	0.85
1:A:206:TYR:CE2	1:A:431:ARG:NH1	2.46	0.83
1:A:69:LEU:HD12	1:A:95:ALA:HB3	1.61	0.81
1:D:395:TRP:CZ2	2:D:451:GOL:H32	2.17	0.80
1:B:195:GLU:CD	1:D:282:ARG:HH22	1.92	0.78
1:A:350:LEU:HD23	1:A:434:LEU:CD1	2.14	0.77
1:B:47:THR:HG23	1:B:57:VAL:HG22	1.67	0.76
1:A:380:VAL:HG22	1:A:387:TYR:CD1	2.24	0.73
1:A:47:THR:HG23	1:A:57:VAL:HG22	1.70	0.73
1:A:69:LEU:HD12	1:A:95:ALA:CB	2.18	0.72
1:C:350:LEU:HD23	1:C:434:LEU:CD1	2.20	0.71
1:D:350:LEU:HD23	1:D:434:LEU:CD1	2.20	0.70
1:C:12:VAL:O	1:C:12:VAL:CG1	2.41	0.69
1:C:36:GLU:HG3	1:C:110:LEU:CD2	2.23	0.69
1:C:375:LEU:O	1:C:376:MSE:CG	2.41	0.68
1:B:326:TYR:O	1:B:341:THR:HG21	1.94	0.67
1:C:36:GLU:HG3	1:C:110:LEU:HD21	1.77	0.66
1:B:350:LEU:HD23	1:B:434:LEU:HD12	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:69:LEU:CD1	1:D:95:ALA:HB3	2.28	0.64
1:D:239:GLN:H	2:D:451:GOL:H2	1.61	0.64
1:A:69:LEU:CD1	1:A:95:ALA:HB3	2.26	0.64
1:C:105:ASP:HB3	1:C:108:ALA:HB2	1.79	0.64
1:B:417:LEU:C	1:B:417:LEU:HD23	2.22	0.64
1:B:8:VAL:HG22	1:B:41:LEU:HD21	1.80	0.64
1:B:350:LEU:HD23	1:B:434:LEU:HD11	1.80	0.63
1:A:350:LEU:HD23	1:A:434:LEU:HD12	1.79	0.63
1:B:8:VAL:HG22	1:B:41:LEU:CD2	2.29	0.62
1:C:256:PHE:HB2	1:C:434:LEU:HD21	1.81	0.62
1:A:13:LEU:HB2	1:A:14:PRO:HD3	1.80	0.62
1:C:8:VAL:HG13	1:C:41:LEU:HD23	1.80	0.62
1:C:375:LEU:C	1:C:376:MSE:CG	2.51	0.61
1:D:69:LEU:HD12	1:D:95:ALA:HB3	1.82	0.61
1:A:429:GLU:OE2	1:A:431:ARG:NE	2.35	0.60
1:B:415:ARG:HD3	3:B:495:HOH:O	2.00	0.60
1:C:417:LEU:C	1:C:417:LEU:HD23	2.25	0.60
1:A:206:TYR:CZ	1:A:431:ARG:NH1	2.69	0.60
1:A:301:GLY:O	1:A:348:HIS:HB3	2.02	0.60
1:A:166:GLN:HG3	1:A:184:THR:HG22	1.84	0.60
1:A:375:LEU:O	1:A:376:MSE:HG3	2.00	0.60
1:C:350:LEU:HD23	1:C:434:LEU:HD12	1.83	0.60
1:B:105:ASP:HB3	1:B:108:ALA:HB2	1.82	0.59
1:A:239:GLN:HA	1:A:239:GLN:OE1	2.01	0.59
1:A:243:VAL:HB	1:A:244:PRO:HD2	1.83	0.59
1:D:417:LEU:C	1:D:417:LEU:HD23	2.26	0.59
1:A:164:THR:HG1	1:A:186:HIS:CE1	2.19	0.59
1:A:417:LEU:HD23	1:A:417:LEU:C	2.28	0.59
1:C:301:GLY:O	1:C:348:HIS:HB3	2.02	0.59
1:A:134:GLU:CD	1:B:364:ARG:HD3	2.28	0.59
1:D:350:LEU:HD23	1:D:434:LEU:HD11	1.83	0.59
1:B:162:LEU:HD12	1:B:188:LEU:HD13	1.83	0.58
1:B:200:MSE:CE	1:D:419:ARG:NH2	2.66	0.58
1:A:9:GLY:O	1:A:12:VAL:HG22	2.04	0.58
1:B:301:GLY:O	1:B:348:HIS:HB3	2.03	0.58
1:D:69:LEU:HD12	1:D:95:ALA:CB	2.33	0.58
1:A:162:LEU:HD12	1:A:188:LEU:HD13	1.86	0.57
1:D:105:ASP:HB3	1:D:108:ALA:HB2	1.85	0.57
1:A:243:VAL:HB	1:A:244:PRO:CD	2.35	0.57
1:C:251:ILE:HG23	1:C:375:LEU:HD12	1.87	0.57
1:B:195:GLU:OE2	1:D:282:ARG:NH1	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:427:ASP:OD1	1:D:427:ASP:N	2.38	0.56
1:C:12:VAL:O	1:C:12:VAL:HG13	2.04	0.56
1:A:10:PRO:HA	1:A:13:LEU:HG	1.88	0.56
1:A:35:CYS:CB	1:A:69:LEU:HD23	2.36	0.56
1:B:200:MSE:HE3	1:D:419:ARG:NH2	2.21	0.56
1:B:256:PHE:HB2	1:B:434:LEU:HD21	1.86	0.56
1:D:232:ARG:HG2	3:D:494:HOH:O	2.05	0.56
1:B:423:ALA:HA	2:B:451:GOL:H31	1.88	0.55
1:A:105:ASP:HB3	1:A:108:ALA:HB2	1.89	0.55
1:D:336:LEU:HB3	1:D:338:LEU:HD12	1.89	0.55
1:D:26:GLU:O	1:D:30:GLN:HG3	2.07	0.55
1:A:256:PHE:HB2	1:A:434:LEU:HD21	1.88	0.54
1:B:423:ALA:HB1	2:B:451:GOL:H32	1.89	0.54
1:A:26:GLU:O	1:A:30:GLN:HG3	2.07	0.54
1:A:366:MSE:HG3	1:A:440:LEU:HD11	1.89	0.54
1:D:173:ASN:HB3	1:D:175:HIS:H	1.73	0.53
1:A:35:CYS:HB2	1:A:69:LEU:HD23	1.89	0.53
1:C:350:LEU:HD23	1:C:434:LEU:HD11	1.87	0.53
1:B:8:VAL:HG13	1:B:41:LEU:HD23	1.91	0.53
1:B:265:GLY:HA3	1:B:332:SER:OG	2.09	0.53
1:A:196:GLY:HA3	1:A:233:LEU:HD13	1.90	0.53
1:A:134:GLU:OE2	1:B:364:ARG:HD3	2.10	0.52
1:A:84:GLY:O	1:A:88:ILE:HG13	2.08	0.52
1:A:425:GLY:O	1:A:437:GLY:HA3	2.09	0.52
1:C:172:GLY:HA2	1:C:177:PHE:O	2.10	0.52
1:B:111:ALA:HB3	1:B:116:MSE:HE1	1.91	0.52
1:D:173:ASN:HD22	1:D:175:HIS:HB2	1.75	0.52
1:B:124:ASP:HB3	1:B:297:VAL:HG13	1.92	0.52
1:D:35:CYS:H	1:D:174:SER:HB3	1.74	0.52
1:B:383:VAL:HG11	1:C:200:MSE:HE1	1.92	0.51
1:C:116:MSE:HE3	1:C:175:HIS:HD2	1.75	0.51
1:B:350:LEU:HB3	1:B:434:LEU:HD12	1.93	0.51
1:B:380:VAL:HG22	1:B:387:TYR:CD1	2.45	0.51
1:C:166:GLN:HG3	1:C:184:THR:HG22	1.90	0.51
1:D:256:PHE:HB2	1:D:434:LEU:HD21	1.91	0.51
1:D:124:ASP:HB3	1:D:297:VAL:HG13	1.92	0.51
1:B:206:TYR:CE2	1:B:431:ARG:NH1	2.79	0.51
1:A:253:VAL:HG13	1:A:434:LEU:HD23	1.92	0.51
1:C:336:LEU:HB3	1:C:338:LEU:HD12	1.92	0.51
1:C:188:LEU:HD23	1:C:221:ILE:HD12	1.92	0.50
1:D:164:THR:HG23	1:D:164:THR:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:64:GLY:HA2	1:D:81:SER:HA	1.92	0.50
1:D:195:GLU:HG2	1:D:200:MSE:HG3	1.93	0.50
1:C:204:TYR:CD1	1:C:204:TYR:C	2.89	0.50
1:D:350:LEU:HB3	1:D:434:LEU:HD12	1.93	0.50
1:A:204:TYR:CD2	1:A:204:TYR:N	2.76	0.50
1:B:36:GLU:HG2	1:B:171:TYR:CE1	2.47	0.50
1:B:94:ALA:O	1:B:95:ALA:C	2.55	0.50
1:D:188:LEU:HD23	1:D:221:ILE:HD12	1.92	0.50
1:A:299:ALA:HA	3:A:457:HOH:O	2.11	0.50
1:D:47:THR:HG23	1:D:57:VAL:HG22	1.94	0.50
1:C:45:LEU:HD12	1:C:58:GLU:O	2.12	0.49
1:D:45:LEU:HD12	1:D:58:GLU:O	2.13	0.49
1:A:158:ASP:OD2	1:A:202:ARG:NH1	2.46	0.49
1:A:8:VAL:HG13	1:A:41:LEU:CD2	2.28	0.49
1:B:204:TYR:CD1	1:B:204:TYR:C	2.90	0.49
1:D:253:VAL:HG13	1:D:434:LEU:HD23	1.95	0.49
1:A:379:GLY:HA3	1:A:388:SER:O	2.13	0.48
1:D:196:GLY:HA3	1:D:233:LEU:HD13	1.94	0.48
1:A:199:GLN:OE1	1:A:234:GLY:N	2.40	0.48
1:C:448:GLY:O	1:C:449:ARG:CB	2.61	0.48
1:D:204:TYR:CD1	1:D:204:TYR:C	2.91	0.48
1:D:395:TRP:CH2	2:D:451:GOL:H32	2.48	0.48
1:B:37:VAL:HG12	1:B:38:ALA:N	2.29	0.48
1:D:36:GLU:HG2	1:D:171:TYR:CE1	2.49	0.48
1:B:154:VAL:HG21	1:B:226:ALA:HB1	1.96	0.48
1:A:275:LEU:HD23	1:A:275:LEU:HA	1.64	0.48
1:B:200:MSE:HE1	1:D:419:ARG:NH2	2.29	0.48
1:D:12:VAL:HG13	1:D:12:VAL:O	2.12	0.48
1:D:196:GLY:O	1:D:199:GLN:N	2.41	0.48
1:A:42:GLU:HG2	1:A:43:GLN:N	2.25	0.47
1:A:67:ILE:HD11	1:A:69:LEU:HD21	1.95	0.47
1:C:253:VAL:HG13	1:C:434:LEU:HD23	1.96	0.47
1:C:54:VAL:O	1:C:54:VAL:HG12	2.06	0.47
1:C:328:LEU:HD22	1:C:332:SER:HB3	1.96	0.47
1:A:435:HIS:CD2	1:A:435:HIS:N	2.83	0.47
1:B:363:ILE:HA	1:B:366:MSE:CE	2.44	0.47
1:C:253:VAL:HA	1:C:434:LEU:HD23	1.97	0.47
1:B:423:ALA:CB	2:B:451:GOL:H32	2.43	0.47
1:B:428:ILE:HD12	1:B:428:ILE:N	2.30	0.47
1:D:248:ALA:O	1:D:249:PRO:C	2.57	0.47
1:A:375:LEU:C	1:A:376:MSE:HG3	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:ASP:HB3	1:C:297:VAL:HG13	1.97	0.47
1:D:286:GLU:O	1:D:320:ASN:HA	2.15	0.47
1:A:124:ASP:HB3	1:A:297:VAL:HG13	1.97	0.47
1:B:155:THR:C	1:B:156:LYS:HG3	2.39	0.47
1:C:55:GLU:C	1:C:56:THR:CG2	2.88	0.47
1:C:84:GLY:O	1:C:88:ILE:HG13	2.15	0.47
1:C:366:MSE:HE3	1:C:366:MSE:HB3	1.81	0.47
2:C:451:GOL:H31	3:C:482:HOH:O	2.15	0.47
1:B:45:LEU:HD12	1:B:58:GLU:O	2.15	0.47
1:D:388:SER:O	1:D:389:ARG:HG2	2.14	0.47
1:B:36:GLU:HG2	1:B:171:TYR:HE1	1.80	0.46
1:A:177:PHE:O	1:A:178:ILE:HD13	2.15	0.46
1:D:380:VAL:HG22	1:D:387:TYR:CD1	2.50	0.46
1:B:177:PHE:O	1:B:178:ILE:HD13	2.15	0.46
1:A:111:ALA:HB3	1:A:116:MSE:HE1	1.98	0.46
1:D:166:GLN:HG3	1:D:184:THR:HG22	1.98	0.46
1:D:206:TYR:CE2	1:D:431:ARG:NH1	2.84	0.46
1:D:301:GLY:O	1:D:348:HIS:HB3	2.15	0.46
1:C:67:ILE:HD11	1:C:69:LEU:HD21	1.98	0.46
1:D:278:ALA:O	1:D:279:LEU:C	2.58	0.46
1:B:328:LEU:HD22	1:B:332:SER:HB3	1.98	0.45
1:A:429:GLU:OE1	1:A:429:GLU:HA	2.17	0.45
1:A:278:ALA:O	1:A:279:LEU:C	2.58	0.45
1:D:238:VAL:HB	2:D:451:GOL:H31	1.99	0.45
1:D:282:ARG:HB3	1:D:419:ARG:NH1	2.31	0.45
1:A:170:VAL:HG22	1:A:180:GLY:HA3	1.99	0.45
1:C:170:VAL:HG22	1:C:180:GLY:HA3	1.99	0.45
1:C:55:GLU:C	1:C:56:THR:HG23	2.40	0.45
1:D:69:LEU:HD13	1:D:96:LEU:HG	1.99	0.45
1:A:173:ASN:OD1	1:A:175:HIS:HB2	2.16	0.45
1:B:336:LEU:HB3	1:B:338:LEU:HD12	1.98	0.45
1:B:200:MSE:SE	1:C:381:ASN:HD21	2.50	0.44
1:A:36:GLU:HG2	1:A:171:TYR:CE1	2.53	0.44
1:A:350:LEU:HD23	1:A:434:LEU:HD11	1.98	0.44
1:A:13:LEU:N	1:A:14:PRO:CD	2.81	0.44
1:B:156:LYS:HB2	1:B:193:ILE:HB	2.00	0.44
1:C:265:GLY:HA3	1:C:332:SER:OG	2.18	0.44
1:C:380:VAL:HG22	1:C:387:TYR:CD1	2.53	0.44
1:A:20:VAL:HG21	1:A:170:VAL:HG12	2.00	0.43
1:A:350:LEU:HB3	1:A:434:LEU:HD12	1.99	0.43
1:B:278:ALA:C	1:B:323:LEU:HD23	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:363:ILE:HA	1:B:366:MSE:HE2	2.00	0.43
1:B:8:VAL:O	1:B:166:GLN:OE1	2.36	0.43
1:B:69:LEU:CD1	1:B:96:LEU:HG	2.48	0.43
1:D:29:ARG:NH2	3:D:487:HOH:O	2.51	0.43
1:B:161:THR:O	1:B:188:LEU:HD12	2.18	0.43
1:B:243:VAL:HB	1:B:244:PRO:HD2	2.01	0.43
1:C:371:LEU:HD12	1:C:372:VAL:N	2.34	0.43
1:D:170:VAL:HG22	1:D:180:GLY:HA3	2.01	0.43
1:D:350:LEU:HD23	1:D:434:LEU:HD12	1.97	0.43
1:A:265:GLY:HA3	1:A:332:SER:OG	2.19	0.43
1:A:239:GLN:OE1	1:A:239:GLN:CA	2.62	0.43
1:C:111:ALA:HB3	1:C:116:MSE:HE1	2.01	0.43
1:C:394:TYR:CD2	1:C:394:TYR:N	2.87	0.43
1:A:188:LEU:HD23	1:A:221:ILE:HD12	2.01	0.42
1:A:35:CYS:HB3	1:A:69:LEU:HD23	2.00	0.42
1:B:429:GLU:OE2	1:B:429:GLU:HA	2.19	0.42
1:C:275:LEU:HD23	1:C:275:LEU:HA	1.77	0.42
1:B:170:VAL:HG22	1:B:180:GLY:HA3	2.00	0.42
1:C:60:ASN:HB2	3:C:523:HOH:O	2.20	0.42
1:D:256:PHE:CE2	1:D:352:VAL:HG21	2.54	0.42
1:A:70:TYR:HA	1:A:74:ARG:O	2.19	0.42
1:A:278:ALA:C	1:A:323:LEU:HD23	2.44	0.42
1:D:328:LEU:HD22	1:D:332:SER:HB3	2.01	0.42
1:B:67:ILE:HD11	1:B:69:LEU:HD21	2.01	0.42
1:B:371:LEU:HD12	1:B:372:VAL:N	2.35	0.42
1:B:279:LEU:HD12	1:B:279:LEU:HA	1.84	0.42
1:A:358:ASP:O	1:A:359:GLN:C	2.62	0.42
1:B:417:LEU:C	1:B:417:LEU:CD2	2.92	0.42
1:C:417:LEU:C	1:C:417:LEU:CD2	2.92	0.41
1:D:275:LEU:HA	1:D:275:LEU:HD23	1.88	0.41
1:A:8:VAL:HG22	1:A:41:LEU:HD21	2.02	0.41
1:A:286:GLU:O	1:A:320:ASN:HA	2.20	0.41
1:B:417:LEU:HD23	1:B:417:LEU:O	2.19	0.41
1:C:319:GLU:O	1:C:320:ASN:C	2.61	0.41
1:B:243:VAL:HB	1:B:244:PRO:CD	2.50	0.41
1:D:13:LEU:HB2	1:D:14:PRO:HD3	2.03	0.41
1:D:19:GLN:O	1:D:23:ILE:HG13	2.20	0.41
1:B:13:LEU:HB2	1:B:14:PRO:HD3	2.02	0.41
1:D:395:TRP:CH2	2:D:451:GOL:C3	3.03	0.41
1:A:12:VAL:O	1:A:13:LEU:C	2.63	0.41
1:D:363:ILE:HA	1:D:366:MSE:HE2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:GLY:O	1:A:389:ARG:NH2	2.53	0.41
1:A:65:PHE:CD1	1:A:65:PHE:C	2.99	0.41
1:C:206:TYR:CE2	1:C:431:ARG:NH1	2.89	0.41
1:C:256:PHE:CE2	1:C:352:VAL:HG21	2.56	0.41
1:A:137:VAL:O	1:A:138:GLU:C	2.64	0.40
1:A:190:CYS:O	1:A:204:TYR:CB	2.69	0.40
1:B:200:MSE:CE	1:D:419:ARG:HH22	2.34	0.40
1:B:253:VAL:HG13	1:B:434:LEU:HD23	2.02	0.40
1:B:286:GLU:O	1:B:320:ASN:HA	2.21	0.40
1:D:36:GLU:HG2	1:D:171:TYR:HE1	1.84	0.40
1:B:364:ARG:NH2	3:B:558:HOH:O	2.55	0.40
1:B:429:GLU:OE1	1:B:431:ARG:NH2	2.54	0.40
1:D:127:HIS:C	1:D:184:THR:HG21	2.47	0.40
1:D:265:GLY:HA3	1:D:332:SER:OG	2.22	0.40
1:B:200:MSE:HE1	1:D:419:ARG:HH22	1.85	0.40
1:D:111:ALA:HB3	1:D:116:MSE:HE1	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	440/449 (98%)	434 (99%)	6 (1%)	0	100	100
1	B	439/449 (98%)	434 (99%)	5 (1%)	0	100	100
1	C	440/449 (98%)	437 (99%)	3 (1%)	0	100	100
1	D	439/449 (98%)	434 (99%)	5 (1%)	0	100	100
All	All	1758/1796 (98%)	1739 (99%)	19 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	330/331 (100%)	327 (99%)	3 (1%)	70	87
1	B	332/331 (100%)	330 (99%)	2 (1%)	78	91
1	C	329/331 (99%)	323 (98%)	6 (2%)	51	78
1	D	331/331 (100%)	328 (99%)	3 (1%)	70	87
All	All	1322/1324 (100%)	1308 (99%)	14 (1%)	65	85

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	VAL
1	A	376	MSE
1	A	382	LEU
1	B	69	LEU
1	B	122	GLU
1	C	12	VAL
1	C	69	LEU
1	C	173	ASN
1	C	195	GLU
1	C	197	GLU
1	C	371	LEU
1	D	12	VAL
1	D	33	SER
1	D	67	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	165	HIS
1	A	402	GLN
1	B	73	GLN
1	C	175	HIS

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Mol	Chain	Res	Type
1	C	281	GLN
1	D	43	GLN
1	D	51	GLN
1	D	173	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](#)

9 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	GOL	C	450	-	5,5,5	0.65	0	5,5,5	0.30	0
2	GOL	B	450	-	5,5,5	1.13	1 (20%)	5,5,5	1.07	0
2	GOL	D	451	-	5,5,5	0.87	0	5,5,5	1.26	0
2	GOL	D	452	-	5,5,5	0.71	0	5,5,5	0.47	0
2	GOL	A	450	-	5,5,5	1.15	1 (20%)	5,5,5	1.18	1 (20%)
2	GOL	C	451	-	5,5,5	0.70	0	5,5,5	0.24	0
2	GOL	D	450	-	5,5,5	0.23	0	5,5,5	0.99	0
2	GOL	B	451	-	5,5,5	0.43	0	5,5,5	1.62	1 (20%)
2	GOL	A	451	-	5,5,5	0.91	0	5,5,5	0.60	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	GOL	C	450	-	-	0/4/4/4	-
2	GOL	B	450	-	-	4/4/4/4	-
2	GOL	D	451	-	-	3/4/4/4	-
2	GOL	D	452	-	-	1/4/4/4	-
2	GOL	A	450	-	-	3/4/4/4	-
2	GOL	C	451	-	-	2/4/4/4	-
2	GOL	D	450	-	-	1/4/4/4	-
2	GOL	B	451	-	-	0/4/4/4	-
2	GOL	A	451	-	-	4/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	450	GOL	O2-C2	-2.30	1.36	1.43
2	B	450	GOL	O2-C2	-2.21	1.36	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	451	GOL	O3-C3-C2	2.53	121.77	110.38
2	A	450	GOL	O2-C2-C1	-2.45	99.03	109.18

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	451	GOL	C1-C2-C3-O3
2	A	451	GOL	O2-C2-C3-O3
2	B	450	GOL	O1-C1-C2-C3
2	C	451	GOL	O1-C1-C2-C3
2	D	451	GOL	C1-C2-C3-O3
2	B	450	GOL	O1-C1-C2-O2
2	A	450	GOL	C1-C2-C3-O3
2	A	451	GOL	O1-C1-C2-C3
2	C	451	GOL	O1-C1-C2-O2
2	A	451	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	D	451	GOL	O2-C2-C3-O3
2	D	450	GOL	O1-C1-C2-O2
2	D	451	GOL	O1-C1-C2-O2
2	A	450	GOL	O1-C1-C2-C3
2	B	450	GOL	C1-C2-C3-O3
2	A	450	GOL	O1-C1-C2-O2
2	B	450	GOL	O2-C2-C3-O3
2	D	452	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	451	GOL	5	0
2	C	451	GOL	1	0
2	B	451	GOL	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	435/449 (96%)	-0.07	3 (0%) 84 83	13, 27, 50, 80	0
1	B	434/449 (96%)	-0.08	2 (0%) 87 86	12, 26, 49, 61	0
1	C	435/449 (96%)	0.05	7 (1%) 70 68	17, 32, 53, 74	0
1	D	434/449 (96%)	0.05	5 (1%) 76 75	18, 33, 54, 71	0
All	All	1738/1796 (96%)	-0.01	17 (0%) 79 78	12, 29, 52, 80	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	449	ARG	5.2
1	B	35	CYS	4.0
1	C	449	ARG	4.0
1	A	130	SER	3.6
1	C	35	CYS	3.2
1	D	181	TYR	2.9
1	C	53	GLU	2.9
1	B	60	ASN	2.9
1	D	8	VAL	2.6
1	A	204	TYR	2.5
1	D	139	ARG	2.4
1	C	59	PHE	2.3
1	D	202	ARG	2.2
1	C	169	ARG	2.0
1	C	415	ARG	2.0
1	D	53	GLU	2.0
1	C	181	TYR	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	GOL	C	451	6/6	0.68	0.19	44,58,69,70	0
2	GOL	D	451	6/6	0.74	0.18	53,56,60,61	0
2	GOL	D	450	6/6	0.77	0.16	34,40,48,58	0
2	GOL	A	451	6/6	0.77	0.19	33,44,47,59	0
2	GOL	D	452	6/6	0.78	0.18	40,55,57,60	0
2	GOL	A	450	6/6	0.79	0.16	33,40,45,52	0
2	GOL	C	450	6/6	0.80	0.12	29,39,44,47	0
2	GOL	B	450	6/6	0.91	0.12	34,41,50,51	0
2	GOL	B	451	6/6	0.92	0.10	39,43,50,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.