



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:39 PM UTC

PDB ID : 1WOG / pdb_00001wog
Title : Crystal Structure of Agmatinase Reveals Structural Conservation and Inhibition Mechanism of the Ureohydrolase Superfamily
Authors : Ahn, H.J.; Kim, K.H.; Lee, J.; Ha, J.-Y.; Lee, H.H.; Kim, D.; Yoon, H.-J.; Kwon, A.-R.; Suh, S.W.
Deposited on : 2004-08-18
Resolution : 1.80 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

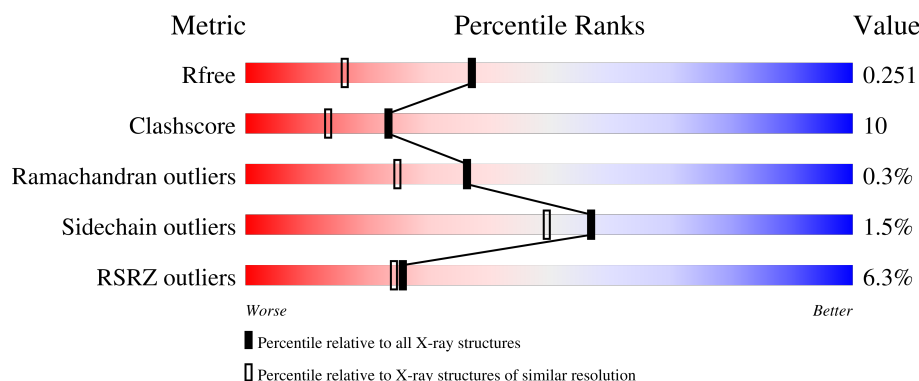
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	305	4% 79% 19% ..
1	B	305	6% 78% 19% ..
1	C	305	5% 80% 19% ..
1	D	305	9% 82% 17% ..
1	E	305	8% 81% 17% ..

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Mol	Chain	Length	Quality of chain
1	F	305	

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
3	16D	A	1401	-	-	X	-
3	16D	B	1402	-	-	X	-
3	16D	C	1403	-	-	X	-
3	16D	D	1404	-	-	X	-
3	16D	E	1405	-	-	X	-
3	16D	F	1406	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14625 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 agmatinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	0	0	0
			2287	1446	409	425	7			
1	B	303	Total	C	N	O	S	0	0	0
			2287	1446	409	425	7			
1	C	303	Total	C	N	O	S	0	0	0
			2287	1446	409	425	7			
1	D	303	Total	C	N	O	S	0	0	0
			2287	1446	409	425	7			
1	E	303	Total	C	N	O	S	0	0	0
			2287	1446	409	425	7			
1	F	303	Total	C	N	O	S	0	0	0
			2287	1446	409	425	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	305	LEU	-	cloning artifact	UNP Q9RZ04
B	305	LEU	-	cloning artifact	UNP Q9RZ04
C	305	LEU	-	cloning artifact	UNP Q9RZ04
D	305	LEU	-	cloning artifact	UNP Q9RZ04
E	305	LEU	-	cloning artifact	UNP Q9RZ04
F	305	LEU	-	cloning artifact	UNP Q9RZ04

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

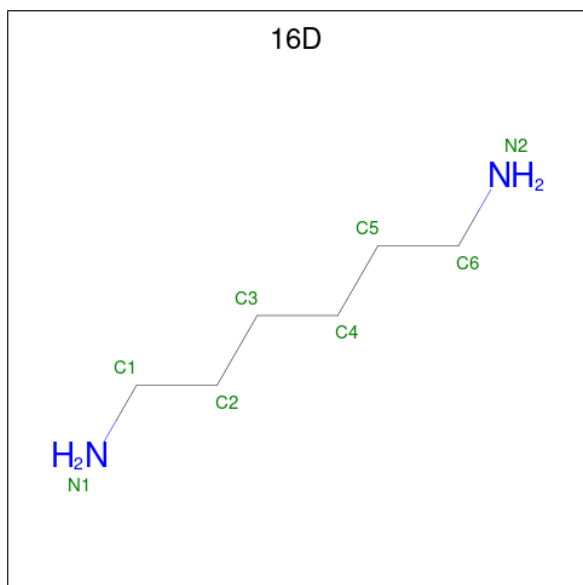
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mn	0	0
			2	2		
2	B	2	Total	Mn	0	0
			2	2		
2	C	2	Total	Mn	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	2	Total	Mn	0	0
			2	2		
2	E	2	Total	Mn	0	0
			2	2		
2	F	2	Total	Mn	0	0
			2	2		

- Molecule 3 is HEXANE-1,6-DIAMINE (CCD ID: 16D) (formula: $C_6H_{16}N_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			8	6	2		
3	B	1	Total	C	N	0	0
			8	6	2		
3	C	1	Total	C	N	0	0
			8	6	2		
3	D	1	Total	C	N	0	0
			8	6	2		
3	E	1	Total	C	N	0	0
			8	6	2		
3	F	1	Total	C	N	0	0
			8	6	2		

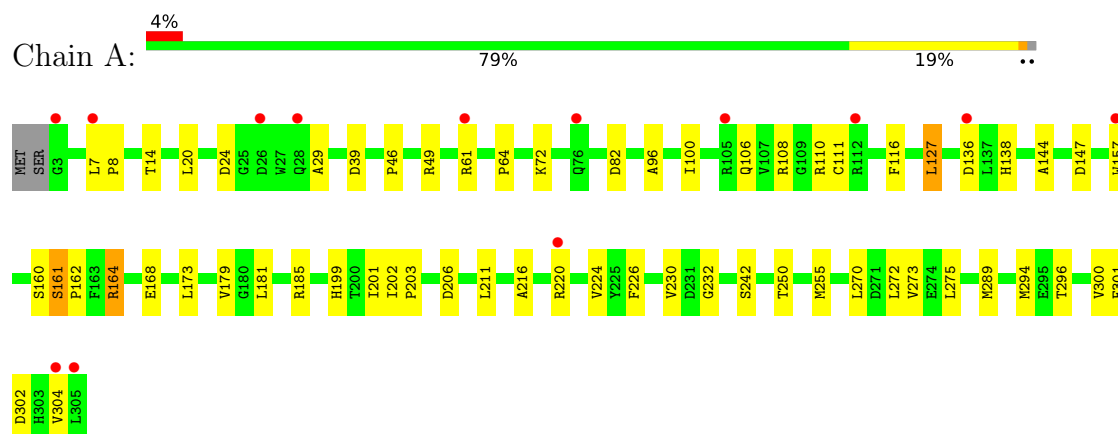
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	169	Total 169	O 169	0	0
4	B	126	Total 126	O 126	0	0
4	C	151	Total 151	O 151	0	0
4	D	113	Total 113	O 113	0	0
4	E	133	Total 133	O 133	0	0
4	F	151	Total 151	O 151	0	0

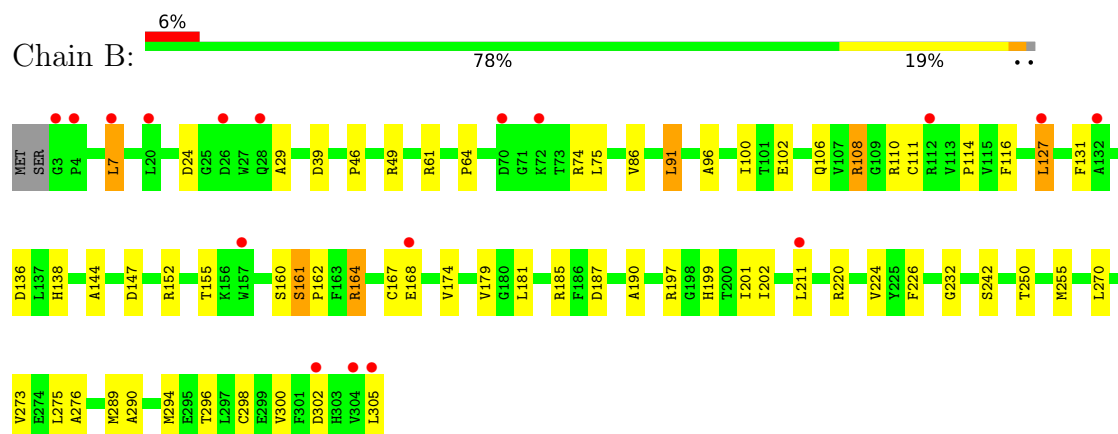
3 Residue-property plots [i](#)

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.

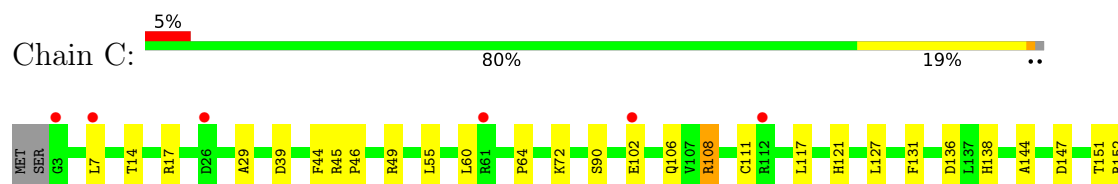
• Molecule 1: agmatinase



• Molecule 1: agmatinase

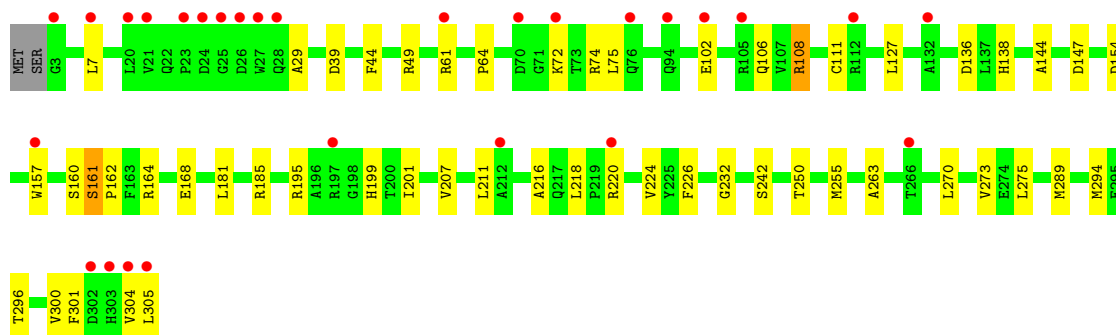
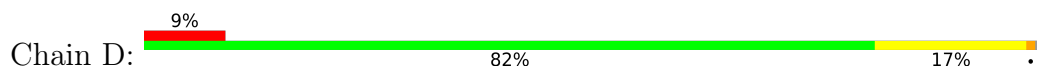


• Molecule 1: agmatinase

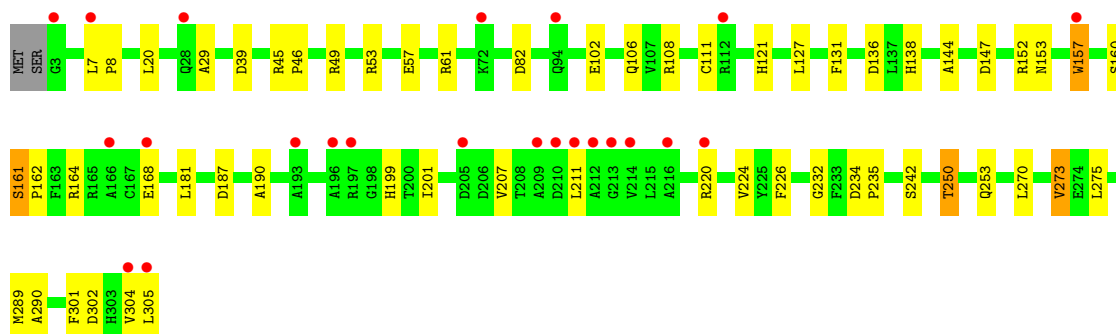
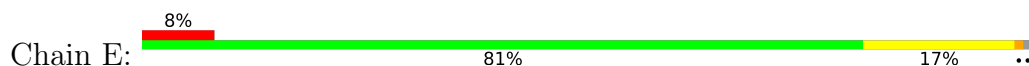




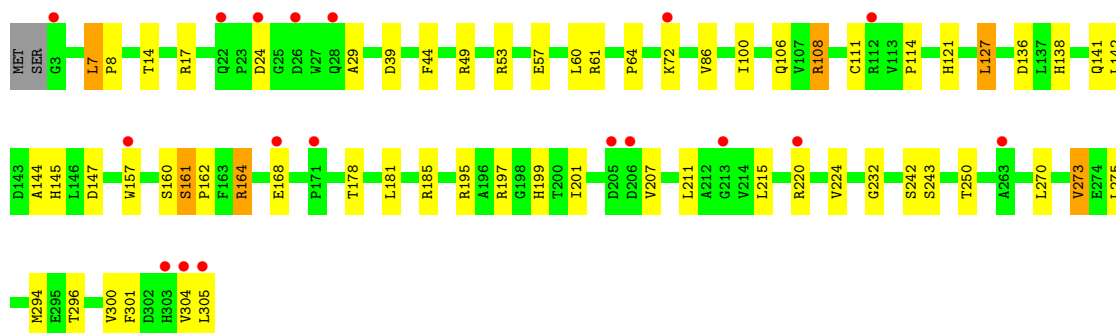
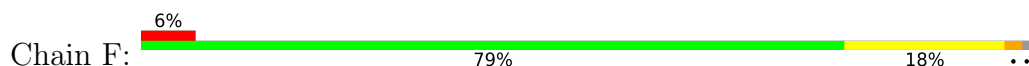
● Molecule 1: agmatinase



● Molecule 1: agmatinase



● Molecule 1: agmatinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.77Å 131.44Å 168.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	16.46 – 1.80 16.46 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.7 (16.46-1.80) 97.7 (16.46-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.62 (at 1.80Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.251 0.224 , 0.251	Depositor DCC
R_{free} test set	16341 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	16.2	Xtriage
Anisotropy	0.236	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.53 , 63.7	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	0.92	EDS
Total number of atoms	14625	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 3.59% 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, 0.333 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: 16D, MN

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	0.39	0/2342	0.92	10/3202 (0.3%)
1	B	0.36	0/2342	0.91	14/3202 (0.4%)
1	C	0.37	0/2342	0.92	13/3202 (0.4%)
1	D	0.35	0/2342	0.92	10/3202 (0.3%)
1	E	0.36	0/2342	0.92	9/3202 (0.3%)
1	F	0.36	0/2342	0.91	12/3202 (0.4%)
All	All	0.36	0/14052	0.92	68/19212 (0.4%)

There are no bond length outliers.

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	LEU	N-CA-C	-8.49	100.36	110.13
1	E	250	THR	N-CA-C	-8.46	99.61	110.53
1	B	250	THR	N-CA-C	-8.32	99.80	110.53
1	E	7	LEU	N-CA-C	-8.26	100.63	110.13
1	D	250	THR	N-CA-C	-7.93	100.30	110.53
1	D	7	LEU	N-CA-C	-7.81	101.15	110.13
1	C	7	LEU	N-CA-C	-7.70	101.28	110.13
1	A	250	THR	N-CA-C	-7.64	100.67	110.53
1	F	250	THR	N-CA-C	-7.33	100.65	110.55
1	C	250	THR	N-CA-C	-7.28	101.14	110.53
1	F	7	LEU	N-CA-C	-6.91	99.98	110.07
1	A	242	SER	N-CA-C	6.75	118.29	111.07
1	F	242	SER	N-CA-C	6.56	118.12	110.97
1	C	224	VAL	N-CA-C	6.46	117.78	108.48
1	D	144	ALA	N-CA-C	-6.44	105.25	113.23
1	C	144	ALA	N-CA-C	-6.39	105.30	113.23
1	A	144	ALA	N-CA-C	-6.34	105.36	113.23
1	B	242	SER	N-CA-C	6.33	117.84	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	224	VAL	N-CA-C	6.25	116.86	108.11
1	F	144	ALA	N-CA-C	-6.16	105.60	113.23
1	D	242	SER	N-CA-C	6.08	117.58	111.07
1	E	242	SER	N-CA-C	6.03	118.35	111.11
1	A	232	GLY	N-CA-C	-5.88	105.50	113.37
1	B	7	LEU	N-CA-C	-5.87	101.50	110.07
1	C	273	VAL	N-CA-C	5.81	117.21	108.96
1	E	273	VAL	N-CA-C	5.74	117.11	108.96
1	C	242	SER	N-CA-C	5.71	117.18	111.07
1	D	232	GLY	N-CA-C	-5.68	105.76	113.37
1	C	14	THR	N-CA-C	-5.67	101.36	110.14
1	B	290	ALA	N-CA-C	-5.66	105.19	111.36
1	D	273	VAL	N-CA-C	5.60	116.91	108.96
1	E	144	ALA	N-CA-C	-5.57	106.33	113.23
1	F	86	VAL	N-CA-C	-5.55	102.41	109.58
1	E	232	GLY	N-CA-C	-5.53	105.97	113.37
1	F	142	LEU	N-CA-C	-5.51	100.26	109.24
1	B	164	ARG	N-CA-C	-5.50	105.37	111.36
1	B	91	LEU	CB-CA-C	-5.49	110.22	116.54
1	C	157	TRP	N-CA-C	5.45	119.24	112.59
1	F	14	THR	N-CA-C	-5.44	101.70	110.14
1	C	232	GLY	N-CA-C	-5.44	106.08	113.37
1	B	224	VAL	N-CA-C	5.44	116.31	108.48
1	B	273	VAL	N-CA-C	5.41	116.70	108.80
1	B	144	ALA	N-CA-C	-5.38	106.21	112.89
1	F	157	TRP	N-CA-C	5.38	119.16	112.59
1	D	224	VAL	N-CA-C	5.38	116.23	108.48
1	A	164	ARG	N-CA-C	-5.36	105.36	111.14
1	E	157	TRP	N-CA-C	5.34	119.10	112.59
1	A	224	VAL	N-CA-C	5.33	116.15	108.48
1	D	185	ARG	N-CA-C	5.32	117.58	108.90
1	E	290	ALA	N-CA-C	-5.31	105.57	111.36
1	F	232	GLY	N-CA-C	-5.29	106.28	113.37
1	C	164	ARG	N-CA-C	-5.27	105.45	111.14
1	F	164	ARG	N-CA-C	-5.25	105.47	111.14
1	E	224	VAL	N-CA-C	5.21	115.98	108.48
1	C	234	ASP	CA-C-N	5.18	124.97	119.28
1	C	234	ASP	C-N-CA	5.18	124.97	119.28
1	D	154	ASP	N-CA-C	5.15	118.83	112.24
1	A	14	THR	N-CA-C	-5.13	101.51	109.72
1	D	157	TRP	N-CA-C	5.13	118.84	112.59
1	B	86	VAL	N-CA-C	-5.11	103.00	109.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	273	VAL	N-CA-C	5.08	116.18	108.96
1	B	276	ALA	CA-C-N	5.08	125.11	119.32
1	B	276	ALA	C-N-CA	5.08	125.11	119.32
1	C	90	SER	N-CA-C	-5.07	101.44	109.96
1	B	174	VAL	N-CA-C	5.05	119.53	112.50
1	A	157	TRP	N-CA-C	5.04	118.74	112.59
1	F	273	VAL	N-CA-C	5.03	116.11	108.96
1	B	232	GLY	N-CA-C	-5.02	106.64	113.37

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	2287	0	2266	47	0
1	B	2287	0	2266	53	0
1	C	2287	0	2266	44	0
1	D	2287	0	2266	48	0
1	E	2287	0	2266	50	0
1	F	2287	0	2266	57	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
3	A	8	0	16	7	0
3	B	8	0	16	6	0
3	C	8	0	16	6	0
3	D	8	0	16	7	0
3	E	8	0	16	6	0
3	F	8	0	16	8	0
4	A	169	0	0	2	0
4	B	126	0	0	4	0
4	C	151	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	113	0	0	3	0
4	E	133	0	0	3	0
4	F	151	0	0	7	0
All	All	14625	0	13692	279	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (279) 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:230:VAL:HG12	1:A:289:MET:HE2	1.21	1.16
1:C:272:LEU:HD21	1:C:289:MET:HE2	1.30	1.08
1:A:272:LEU:HD21	1:A:289:MET:HE3	1.42	0.96
1:A:147:ASP:OD1	3:A:1401:16D:H52	1.70	0.92
1:D:147:ASP:OD1	3:D:1404:16D:H52	1.71	0.90
1:E:181:LEU:HD12	1:E:201:ILE:HG23	1.55	0.88
1:C:272:LEU:HD21	1:C:289:MET:CE	2.05	0.87
1:E:147:ASP:OD1	3:E:1405:16D:H52	1.74	0.86
1:B:147:ASP:OD1	3:B:1402:16D:H52	1.76	0.86
1:C:181:LEU:HD12	1:C:201:ILE:HG23	1.58	0.85
1:C:147:ASP:OD1	3:C:1403:16D:H52	1.79	0.82
1:A:216:ALA:HA	1:A:220:ARG:HH22	1.44	0.81
1:A:272:LEU:HD21	1:A:289:MET:CE	2.11	0.79
1:F:147:ASP:OD1	3:F:1406:16D:H52	1.84	0.77
1:A:147:ASP:OD1	3:A:1401:16D:C3	2.35	0.74
1:D:181:LEU:HD12	1:D:201:ILE:HG23	1.69	0.74
1:D:147:ASP:OD1	3:D:1404:16D:C3	2.36	0.74
1:C:216:ALA:HA	1:C:220:ARG:HH22	1.53	0.73
1:B:147:ASP:OD1	3:B:1402:16D:C3	2.37	0.72
1:E:147:ASP:OD1	3:E:1405:16D:C3	2.38	0.71
1:D:147:ASP:OD1	3:D:1404:16D:H31	1.91	0.71
1:F:207:VAL:HG13	1:F:211:LEU:HD23	1.74	0.69
1:D:61:ARG:HE	1:D:61:ARG:HA	1.57	0.69
1:F:147:ASP:OD1	3:F:1406:16D:C3	2.41	0.68
1:B:147:ASP:OD1	3:B:1402:16D:H31	1.94	0.68
1:B:181:LEU:HD12	1:B:201:ILE:HG23	1.75	0.67
1:A:147:ASP:OD1	3:A:1401:16D:H31	1.95	0.66
1:B:39:ASP:OD2	1:B:49:ARG:HD2	1.95	0.65
1:F:147:ASP:OD1	3:F:1406:16D:H31	1.96	0.65
1:E:181:LEU:CD1	1:E:201:ILE:HG23	2.25	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:PRO:HG3	1:E:49:ARG:NH1	2.11	0.65
1:E:147:ASP:OD1	3:E:1405:16D:H31	1.96	0.65
1:A:300:VAL:O	1:A:304:VAL:HG23	1.97	0.65
1:A:181:LEU:HD12	1:A:201:ILE:HG23	1.78	0.64
1:F:300:VAL:O	1:F:304:VAL:HG23	1.97	0.64
1:F:181:LEU:HD12	1:F:201:ILE:HG23	1.79	0.64
1:D:72:LYS:NZ	1:D:72:LYS:HB3	2.12	0.63
1:C:147:ASP:OD1	3:C:1403:16D:C3	2.47	0.63
1:D:160:SER:HB3	3:D:1404:16D:H12	1.81	0.62
1:A:24:ASP:HB3	1:A:106:GLN:NE2	2.14	0.62
1:D:164:ARG:O	1:D:168:GLU:HG2	2.01	0.61
1:D:301:PHE:HA	1:D:304:VAL:HG23	1.83	0.61
1:A:301:PHE:HA	1:A:304:VAL:HG23	1.81	0.60
1:C:72:LYS:NZ	1:C:72:LYS:HB3	2.15	0.60
1:C:211:LEU:HD13	1:C:211:LEU:C	2.26	0.60
1:A:72:LYS:NZ	1:A:72:LYS:HB3	2.17	0.60
1:F:141:GLN:HG2	1:F:178:THR:HG23	1.82	0.60
1:D:147:ASP:H	3:D:1404:16D:HN11	1.50	0.60
1:F:108:ARG:HG2	1:F:114:PRO:HG3	1.85	0.59
1:A:147:ASP:OD1	3:A:1401:16D:H32	2.01	0.59
1:F:72:LYS:NZ	1:F:72:LYS:HB3	2.17	0.59
1:B:164:ARG:O	1:B:168:GLU:HG2	2.03	0.59
1:D:270:LEU:C	1:D:270:LEU:HD23	2.27	0.58
1:F:270:LEU:HD23	1:F:270:LEU:C	2.28	0.58
1:D:216:ALA:HA	1:D:220:ARG:HH22	1.69	0.58
1:E:39:ASP:CG	1:E:49:ARG:HD2	2.29	0.58
1:E:220:ARG:HG3	1:E:220:ARG:HH11	1.68	0.58
1:B:167:CYS:HB3	1:B:197:ARG:NH1	2.18	0.58
1:C:46:PRO:HG3	1:D:49:ARG:CZ	2.33	0.58
1:C:147:ASP:OD1	3:C:1403:16D:H31	2.04	0.58
1:D:289:MET:HA	1:D:289:MET:HE2	1.85	0.58
1:E:147:ASP:H	3:E:1405:16D:HN11	1.51	0.58
1:E:147:ASP:OD1	3:E:1405:16D:H32	2.04	0.57
1:B:49:ARG:HD3	1:E:46:PRO:HG2	1.86	0.57
1:B:181:LEU:HD12	1:B:201:ILE:CG2	2.33	0.57
1:C:289:MET:HE1	1:C:292:LEU:CD2	2.34	0.57
1:B:147:ASP:OD1	3:B:1402:16D:H32	2.04	0.57
1:A:147:ASP:H	3:A:1401:16D:HN11	1.51	0.57
1:E:108:ARG:HD2	4:E:1648:HOH:O	2.03	0.57
1:C:39:ASP:OD2	1:C:49:ARG:HD2	2.05	0.57
1:B:49:ARG:CZ	1:E:46:PRO:HG3	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:SER:HB3	3:C:1403:16D:H12	1.86	0.56
1:C:289:MET:HE1	1:C:292:LEU:HD22	1.87	0.56
1:C:181:LEU:HD12	1:C:201:ILE:CG2	2.32	0.56
1:E:160:SER:HB3	3:E:1405:16D:H12	1.87	0.56
1:B:46:PRO:HG2	1:E:49:ARG:HD3	1.86	0.56
1:A:29:ALA:O	1:A:111:CYS:HA	2.06	0.56
1:A:46:PRO:HG3	1:F:49:ARG:CZ	2.36	0.56
1:F:164:ARG:O	1:F:168:GLU:HG2	2.06	0.55
1:A:39:ASP:OD2	1:A:49:ARG:HD2	2.07	0.55
1:E:108:ARG:HD3	1:E:131:PHE:CE2	2.42	0.55
1:E:161:SER:N	1:E:162:PRO:HD2	2.21	0.55
1:A:164:ARG:O	1:A:168:GLU:HG2	2.07	0.55
1:B:147:ASP:H	3:B:1402:16D:HN11	1.53	0.55
1:A:216:ALA:HA	1:A:220:ARG:NH2	2.19	0.55
1:E:270:LEU:C	1:E:270:LEU:HD23	2.31	0.55
1:D:161:SER:N	1:D:162:PRO:HD2	2.22	0.55
1:B:298:CYS:O	1:B:302:ASP:HB2	2.06	0.54
1:B:161:SER:N	1:B:162:PRO:HD2	2.23	0.54
1:A:203:PRO:HD2	1:A:206:ASP:OD2	2.08	0.54
1:C:39:ASP:CG	1:C:49:ARG:HD2	2.33	0.54
1:C:301:PHE:HA	1:C:304:VAL:HG23	1.89	0.54
1:A:39:ASP:CG	1:A:49:ARG:HD2	2.32	0.54
1:C:102:GLU:O	1:C:106:GLN:HG3	2.07	0.54
1:F:53:ARG:O	1:F:57:GLU:HG3	2.08	0.54
1:B:289:MET:HE2	1:B:289:MET:HA	1.90	0.54
1:C:185:ARG:HG3	4:C:1756:HOH:O	2.07	0.53
1:D:147:ASP:OD1	3:D:1404:16D:H32	2.07	0.53
1:E:39:ASP:OD2	1:E:49:ARG:HD2	2.08	0.53
1:E:164:ARG:O	1:E:168:GLU:HG2	2.08	0.53
1:F:181:LEU:HD12	1:F:201:ILE:CG2	2.38	0.53
1:F:17:ARG:HG3	1:F:60:LEU:HD22	1.90	0.53
1:A:8:PRO:HG2	1:A:61:ARG:NE	2.24	0.53
1:C:161:SER:N	1:C:162:PRO:HD2	2.23	0.53
1:F:305:LEU:HD13	4:F:1753:HOH:O	2.09	0.53
1:C:181:LEU:CD1	1:C:201:ILE:HG23	2.34	0.53
1:B:46:PRO:HG3	1:E:49:ARG:CZ	2.39	0.53
1:C:270:LEU:HD23	1:C:270:LEU:C	2.34	0.53
1:C:46:PRO:HG2	1:D:49:ARG:HD3	1.90	0.52
1:B:7:LEU:HD23	4:B:1718:HOH:O	2.08	0.52
1:F:136:ASP:O	1:F:138:HIS:HD2	1.91	0.52
1:E:161:SER:N	1:E:162:PRO:CD	2.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:141:GLN:HG3	4:F:1735:HOH:O	2.08	0.52
1:A:270:LEU:HD23	1:A:270:LEU:C	2.35	0.52
1:F:147:ASP:OD1	3:F:1406:16D:H32	2.09	0.52
1:B:211:LEU:HD13	1:B:211:LEU:C	2.35	0.51
1:B:270:LEU:C	1:B:270:LEU:HD23	2.35	0.51
1:B:161:SER:N	1:B:162:PRO:CD	2.73	0.51
1:A:230:VAL:HG12	1:A:289:MET:CE	2.14	0.51
1:D:102:GLU:O	1:D:106:GLN:HG3	2.09	0.51
1:E:187:ASP:HB3	1:E:190:ALA:HB3	1.93	0.51
1:C:108:ARG:HD3	1:C:131:PHE:CE2	2.46	0.51
1:A:199:HIS:HD2	4:A:1682:HOH:O	1.93	0.51
1:B:116:PHE:CB	1:B:127:LEU:HD21	2.40	0.51
1:E:102:GLU:O	1:E:106:GLN:HG3	2.11	0.51
1:C:161:SER:N	1:C:162:PRO:CD	2.74	0.51
1:B:181:LEU:CD1	1:B:201:ILE:HG23	2.41	0.51
1:E:181:LEU:HD12	1:E:201:ILE:CG2	2.35	0.51
1:F:161:SER:N	1:F:162:PRO:CD	2.74	0.51
1:C:17:ARG:HG3	1:C:60:LEU:HD22	1.94	0.50
1:B:108:ARG:HD3	1:B:131:PHE:CE2	2.45	0.50
1:D:39:ASP:CG	1:D:49:ARG:HD2	2.36	0.50
1:B:199:HIS:HE1	4:B:1665:HOH:O	1.93	0.50
1:F:211:LEU:HD13	1:F:211:LEU:O	2.11	0.50
1:A:160:SER:HB3	3:A:1401:16D:H12	1.91	0.50
1:B:46:PRO:CG	1:E:49:ARG:HD3	2.42	0.50
1:E:29:ALA:O	1:E:111:CYS:HA	2.12	0.50
1:D:161:SER:N	1:D:162:PRO:CD	2.74	0.50
1:A:46:PRO:HG3	1:F:49:ARG:NH1	2.27	0.50
1:A:161:SER:N	1:A:162:PRO:CD	2.75	0.50
1:C:147:ASP:H	3:C:1403:16D:HN11	1.59	0.50
1:D:211:LEU:C	1:D:211:LEU:HD13	2.37	0.49
1:D:220:ARG:HG3	1:D:220:ARG:HH11	1.77	0.49
1:B:160:SER:HB3	3:B:1402:16D:H12	1.95	0.49
1:C:301:PHE:HA	1:C:304:VAL:CG2	2.43	0.49
1:B:102:GLU:O	1:B:106:GLN:HG3	2.11	0.49
1:C:147:ASP:OD1	3:C:1403:16D:H32	2.12	0.49
1:A:136:ASP:O	1:A:138:HIS:HD2	1.95	0.49
1:D:304:VAL:C	1:D:305:LEU:HD12	2.38	0.49
1:A:211:LEU:C	1:A:211:LEU:HD13	2.37	0.49
1:F:211:LEU:HD13	1:F:211:LEU:C	2.38	0.48
1:B:296:THR:O	1:B:300:VAL:HG23	2.13	0.48
1:F:147:ASP:H	3:F:1406:16D:HN11	1.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:ARG:NH1	1:E:46:PRO:HG3	2.27	0.48
1:E:226:PHE:O	1:E:270:LEU:HA	2.14	0.48
1:C:64:PRO:HD3	1:C:294:MET:HE2	1.95	0.48
1:C:46:PRO:HG3	1:D:49:ARG:NH1	2.27	0.48
1:E:301:PHE:HA	1:E:304:VAL:HG23	1.96	0.48
1:B:136:ASP:O	1:B:138:HIS:HD2	1.97	0.47
1:D:29:ALA:O	1:D:111:CYS:HA	2.14	0.47
1:B:305:LEU:N	1:B:305:LEU:HD12	2.29	0.47
1:B:211:LEU:HD13	1:B:211:LEU:O	2.13	0.47
1:C:72:LYS:HB3	1:C:72:LYS:HZ3	1.79	0.47
1:E:207:VAL:HG13	1:E:211:LEU:HD23	1.96	0.47
1:F:29:ALA:O	1:F:111:CYS:HA	2.15	0.47
1:E:39:ASP:OD1	1:E:49:ARG:HD2	2.13	0.47
1:B:39:ASP:CG	1:B:49:ARG:HD2	2.40	0.47
1:B:74:ARG:O	1:B:75:LEU:HB2	2.13	0.47
1:C:45:ARG:HA	1:C:46:PRO:HD3	1.81	0.47
1:C:102:GLU:HA	1:C:102:GLU:OE2	2.15	0.47
1:D:44:PHE:CD1	1:F:61:ARG:HD2	2.49	0.47
1:E:61:ARG:HD2	1:F:44:PHE:CD1	2.48	0.47
1:F:39:ASP:OD2	1:F:49:ARG:HD2	2.15	0.47
1:F:296:THR:O	1:F:300:VAL:HG23	2.15	0.47
1:E:45:ARG:HA	1:E:46:PRO:HD3	1.80	0.47
1:F:161:SER:N	1:F:162:PRO:HD2	2.30	0.47
1:D:39:ASP:OD1	1:D:49:ARG:HD2	2.15	0.46
1:F:24:ASP:HB3	1:F:106:GLN:NE2	2.30	0.46
1:F:197:ARG:HD2	4:F:1706:HOH:O	2.13	0.46
1:A:179:VAL:HG22	1:A:202:ILE:HD12	1.97	0.46
1:F:305:LEU:HD12	1:F:305:LEU:N	2.31	0.46
1:A:255:MET:HE1	1:A:296:THR:HA	1.96	0.46
1:B:96:ALA:O	1:B:100:ILE:HG13	2.15	0.46
1:F:301:PHE:HA	1:F:304:VAL:HG23	1.96	0.46
4:E:1646:HOH:O	1:F:185:ARG:HG3	2.14	0.46
1:B:46:PRO:HD2	4:B:1695:HOH:O	2.15	0.46
1:D:207:VAL:HG13	1:D:211:LEU:HD23	1.98	0.46
1:A:181:LEU:HD12	1:A:201:ILE:CG2	2.45	0.46
1:B:29:ALA:O	1:B:111:CYS:HA	2.16	0.46
1:A:220:ARG:HH11	1:A:220:ARG:HG3	1.80	0.46
1:F:8:PRO:HG2	1:F:61:ARG:NE	2.31	0.45
1:F:39:ASP:CG	1:F:49:ARG:HD2	2.41	0.45
1:A:161:SER:N	1:A:162:PRO:HD2	2.30	0.45
1:E:136:ASP:O	1:E:138:HIS:HD2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:74:ARG:O	1:D:75:LEU:HB2	2.16	0.45
1:B:116:PHE:HB3	1:B:127:LEU:HD21	1.97	0.45
1:D:226:PHE:O	1:D:270:LEU:HA	2.17	0.45
1:D:255:MET:HE1	1:D:296:THR:HA	1.99	0.45
1:F:199:HIS:HD2	4:F:1677:HOH:O	1.98	0.45
1:D:64:PRO:HD3	1:D:294:MET:HE2	1.98	0.45
1:A:20:LEU:HD12	1:A:82:ASP:O	2.17	0.45
1:D:181:LEU:CD1	1:D:201:ILE:HG23	2.43	0.45
1:F:304:VAL:HG12	1:F:305:LEU:N	2.32	0.45
1:A:39:ASP:OD2	1:A:49:ARG:CD	2.65	0.44
1:A:46:PRO:HG2	1:F:49:ARG:HD3	1.98	0.44
1:A:185:ARG:HG3	4:A:1669:HOH:O	2.17	0.44
1:D:72:LYS:HB3	1:D:72:LYS:HZ2	1.82	0.44
1:E:61:ARG:HD2	1:F:44:PHE:CE1	2.53	0.44
1:D:108:ARG:HD3	4:D:1688:HOH:O	2.17	0.44
1:A:64:PRO:HD3	1:A:294:MET:HE2	2.00	0.44
1:E:289:MET:HE2	1:E:289:MET:HA	1.98	0.44
1:C:151:THR:HG22	4:C:1726:HOH:O	2.16	0.44
1:E:53:ARG:O	1:E:57:GLU:HG3	2.17	0.44
1:F:160:SER:HB3	3:F:1406:16D:H12	1.99	0.44
1:C:136:ASP:O	1:C:138:HIS:HD2	2.01	0.44
1:B:220:ARG:HG3	1:B:220:ARG:HH11	1.82	0.44
1:C:305:LEU:N	1:C:305:LEU:HD12	2.33	0.44
1:E:250:THR:OG1	1:E:253:GLN:HG3	2.18	0.44
1:B:226:PHE:O	1:B:270:LEU:HA	2.17	0.43
1:C:29:ALA:O	1:C:111:CYS:HA	2.17	0.43
1:A:96:ALA:O	1:A:100:ILE:HG13	2.18	0.43
1:A:301:PHE:HA	1:A:304:VAL:CG2	2.48	0.43
1:B:24:ASP:HB3	1:B:106:GLN:NE2	2.34	0.43
1:B:255:MET:HE1	1:B:296:THR:HA	2.00	0.43
1:F:72:LYS:HB3	1:F:72:LYS:HZ3	1.83	0.43
1:D:181:LEU:HD12	1:D:201:ILE:CG2	2.43	0.43
1:B:61:ARG:HD2	1:C:44:PHE:CE1	2.54	0.42
1:D:102:GLU:OE2	1:D:102:GLU:HA	2.19	0.42
1:B:110:ARG:HH11	1:B:110:ARG:HG3	1.85	0.42
1:D:72:LYS:HB3	1:D:72:LYS:HZ3	1.85	0.42
1:C:211:LEU:HD13	1:C:215:LEU:HG	2.02	0.42
1:D:44:PHE:CE1	1:F:61:ARG:HD2	2.54	0.42
1:F:220:ARG:HD3	4:F:1732:HOH:O	2.18	0.42
1:A:110:ARG:HG3	1:A:110:ARG:HH11	1.85	0.42
1:C:203:PRO:HD2	1:C:206:ASP:OD2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:20:LEU:HD12	1:E:82:ASP:O	2.19	0.42
1:F:145:HIS:HB3	3:F:1406:16D:H32	2.02	0.42
1:F:145:HIS:CB	3:F:1406:16D:H32	2.50	0.42
1:F:181:LEU:CD1	1:F:201:ILE:HG23	2.48	0.42
1:D:195:ARG:HH11	1:D:195:ARG:HG2	1.85	0.42
1:E:61:ARG:NH1	1:F:243:SER:OG	2.52	0.42
1:F:195:ARG:HG2	1:F:195:ARG:HH11	1.85	0.42
1:E:121:HIS:CD2	1:E:273:VAL:HG21	2.55	0.42
1:E:152:ARG:HG2	1:E:153:ASN:OD1	2.20	0.42
1:A:226:PHE:O	1:A:270:LEU:HA	2.20	0.41
1:B:179:VAL:HG22	1:B:202:ILE:HD12	2.02	0.41
1:F:197:ARG:NH2	4:F:1688:HOH:O	2.53	0.41
1:A:147:ASP:CG	3:A:1401:16D:H52	2.43	0.41
1:B:64:PRO:HD3	1:B:294:MET:HE2	2.02	0.41
1:B:91:LEU:CD1	1:E:157:TRP:HB2	2.50	0.41
1:B:152:ARG:O	1:B:155:THR:HG22	2.20	0.41
1:D:305:LEU:HD12	1:D:305:LEU:N	2.35	0.41
1:E:199:HIS:HD2	4:E:1662:HOH:O	2.03	0.41
1:E:220:ARG:HG3	1:E:220:ARG:NH1	2.34	0.41
1:E:211:LEU:HD13	1:E:211:LEU:C	2.45	0.41
1:B:187:ASP:HB3	1:B:190:ALA:HB3	2.02	0.41
1:A:116:PHE:CD1	1:A:127:LEU:HG	2.55	0.41
1:F:64:PRO:HD3	1:F:294:MET:HE2	2.02	0.41
1:F:121:HIS:CD2	1:F:273:VAL:HG21	2.56	0.41
1:C:55:LEU:HD13	1:C:117:LEU:HD22	2.03	0.41
1:F:100:ILE:CG2	1:F:127:LEU:HD13	2.50	0.41
1:D:136:ASP:O	1:D:138:HIS:HD2	2.04	0.41
1:D:160:SER:HA	3:D:1404:16D:H31	2.01	0.41
1:D:199:HIS:HE1	4:D:1692:HOH:O	2.04	0.41
1:D:263:ALA:HB2	4:D:1705:HOH:O	2.21	0.41
1:D:301:PHE:HA	1:D:304:VAL:CG2	2.50	0.41
1:F:211:LEU:HD13	1:F:215:LEU:HG	2.01	0.41
1:B:185:ARG:HG3	4:B:1664:HOH:O	2.20	0.41
1:E:8:PRO:HG2	1:E:61:ARG:NE	2.35	0.41
1:D:218:LEU:O	1:D:220:ARG:NH1	2.54	0.40
1:F:141:GLN:CG	1:F:178:THR:HG23	2.50	0.40
1:B:108:ARG:HG2	1:B:114:PRO:HG3	2.02	0.40
1:C:138:HIS:HE1	4:C:1734:HOH:O	2.04	0.40
1:E:234:ASP:HA	1:E:235:PRO:HD3	1.91	0.40
1:F:7:LEU:HD23	4:F:1696:HOH:O	2.22	0.40
1:A:116:PHE:HB3	1:A:127:LEU:HD21	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:ARG:HG2	1:C:153:ASN:OD1	2.21	0.40
1:E:305:LEU:HD12	1:E:305:LEU:N	2.37	0.40
1:D:296:THR:O	1:D:300:VAL:HG23	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	301/305 (99%)	291 (97%)	9 (3%)	1 (0%)	36	25
1	B	301/305 (99%)	292 (97%)	8 (3%)	1 (0%)	36	25
1	C	301/305 (99%)	291 (97%)	9 (3%)	1 (0%)	36	25
1	D	301/305 (99%)	292 (97%)	8 (3%)	1 (0%)	36	25
1	E	301/305 (99%)	293 (97%)	7 (2%)	1 (0%)	36	25
1	F	301/305 (99%)	291 (97%)	9 (3%)	1 (0%)	36	25
All	All	1806/1830 (99%)	1750 (97%)	50 (3%)	6 (0%)	36	25

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	161	SER
1	E	161	SER
1	A	161	SER
1	B	161	SER
1	C	161	SER
1	D	161	SER

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	240/242 (99%)	235 (98%)	5 (2%)	47	36
1	B	240/242 (99%)	237 (99%)	3 (1%)	61	54
1	C	240/242 (99%)	235 (98%)	5 (2%)	47	36
1	D	240/242 (99%)	237 (99%)	3 (1%)	61	54
1	E	240/242 (99%)	237 (99%)	3 (1%)	61	54
1	F	240/242 (99%)	237 (99%)	3 (1%)	61	54
All	All	1440/1452 (99%)	1418 (98%)	22 (2%)	57	49

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	ARG
1	A	127	LEU
1	A	173	LEU
1	A	275	LEU
1	A	302	ASP
1	B	108	ARG
1	B	127	LEU
1	B	275	LEU
1	C	108	ARG
1	C	121	HIS
1	C	127	LEU
1	C	173	LEU
1	C	275	LEU
1	D	108	ARG
1	D	127	LEU
1	D	275	LEU
1	E	127	LEU
1	E	275	LEU
1	E	302	ASP
1	F	108	ARG
1	F	127	LEU
1	F	275	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	138	HIS
1	A	199	HIS
1	B	76	GLN
1	B	138	HIS
1	B	175	HIS
1	B	199	HIS
1	C	76	GLN
1	C	106	GLN
1	C	138	HIS
1	C	175	HIS
1	C	199	HIS
1	C	264	ASN
1	D	76	GLN
1	D	138	HIS
1	D	199	HIS
1	D	223	ASN
1	E	76	GLN
1	E	138	HIS
1	E	175	HIS
1	E	199	HIS
1	F	76	GLN
1	F	106	GLN
1	F	138	HIS
1	F	199	HIS
1	F	222	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 18 ligands modelled in this entry, 12 are monoatomic - leaving 6 for Mogul analysis.

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	16D	E	1405	-	7,7,7	1.76	1 (14%)	6,6,6	1.04	0
3	16D	F	1406	-	7,7,7	1.76	1 (14%)	6,6,6	1.04	0
3	16D	C	1403	-	7,7,7	1.76	1 (14%)	6,6,6	1.04	0
3	16D	B	1402	-	7,7,7	1.76	1 (14%)	6,6,6	1.04	0
3	16D	D	1404	-	7,7,7	1.77	1 (14%)	6,6,6	1.04	0
3	16D	A	1401	-	7,7,7	1.76	1 (14%)	6,6,6	1.04	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	16D	E	1405	-	-	3/5/5/5	-
3	16D	F	1406	-	-	3/5/5/5	-
3	16D	C	1403	-	-	3/5/5/5	-
3	16D	B	1402	-	-	3/5/5/5	-
3	16D	D	1404	-	-	3/5/5/5	-
3	16D	A	1401	-	-	3/5/5/5	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1404	16D	C5-C6	-4.02	1.32	1.51
3	F	1406	16D	C5-C6	-4.02	1.32	1.51
3	B	1402	16D	C5-C6	-4.01	1.32	1.51
3	C	1403	16D	C5-C6	-4.01	1.32	1.51
3	E	1405	16D	C5-C6	-4.01	1.32	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1401	16D	C5-C6	-4.01	1.32	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1401	16D	C3-C4-C5-C6
3	B	1402	16D	C3-C4-C5-C6
3	C	1403	16D	C3-C4-C5-C6
3	D	1404	16D	C3-C4-C5-C6
3	E	1405	16D	C3-C4-C5-C6
3	F	1406	16D	C3-C4-C5-C6
3	D	1404	16D	C4-C5-C6-N2
3	F	1406	16D	C4-C5-C6-N2
3	A	1401	16D	C4-C5-C6-N2
3	B	1402	16D	C4-C5-C6-N2
3	C	1403	16D	C4-C5-C6-N2
3	E	1405	16D	C4-C5-C6-N2
3	A	1401	16D	N1-C1-C2-C3
3	B	1402	16D	N1-C1-C2-C3
3	C	1403	16D	N1-C1-C2-C3
3	D	1404	16D	N1-C1-C2-C3
3	E	1405	16D	N1-C1-C2-C3
3	F	1406	16D	N1-C1-C2-C3

There are no ring outliers.

6 monomers are involved in 40 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	1405	16D	6	0
3	F	1406	16D	8	0
3	C	1403	16D	6	0
3	B	1402	16D	6	0
3	D	1404	16D	7	0
3	A	1401	16D	7	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	303/305 (99%)	0.08	13 (4%)	40	39	7, 14, 29, 44	0
1	B	303/305 (99%)	0.49	17 (5%)	30	29	9, 19, 35, 73	0
1	C	303/305 (99%)	0.20	16 (5%)	32	31	8, 15, 29, 86	0
1	D	303/305 (99%)	0.55	28 (9%)	14	12	10, 19, 37, 87	0
1	E	303/305 (99%)	0.45	23 (7%)	20	18	10, 18, 32, 62	0
1	F	303/305 (99%)	0.33	18 (5%)	28	27	8, 16, 31, 64	0
All	All	1818/1830 (99%)	0.35	115 (6%)	26	24	7, 17, 34, 87	0

All (115) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	305	LEU	8.4
1	C	304	VAL	7.6
1	C	305	LEU	7.3
1	F	305	LEU	7.1
1	D	304	VAL	7.0
1	D	305	LEU	6.7
1	B	305	LEU	5.7
1	B	304	VAL	5.6
1	C	303	HIS	5.4
1	F	304	VAL	5.4
1	E	304	VAL	4.7
1	A	304	VAL	4.5
1	D	102	GLU	4.0
1	F	112	ARG	3.9
1	A	220	ARG	3.8
1	C	302	ASP	3.6
1	D	3	GLY	3.6
1	A	28	GLN	3.6
1	D	302	ASP	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	102	GLU	3.5
1	F	168	GLU	3.5
1	D	220	ARG	3.5
1	B	3	GLY	3.5
1	A	157	TRP	3.4
1	A	305	LEU	3.4
1	D	72	LYS	3.3
1	E	211	LEU	3.3
1	C	220	ARG	3.3
1	F	220	ARG	3.3
1	E	213	GLY	3.2
1	C	3	GLY	3.1
1	F	3	GLY	3.1
1	E	220	ARG	3.1
1	B	211	LEU	3.0
1	F	24	ASP	3.0
1	F	303	HIS	3.0
1	B	28	GLN	3.0
1	D	157	TRP	3.0
1	E	197	ARG	3.0
1	D	24	ASP	2.9
1	E	216	ALA	2.9
1	B	72	LYS	2.9
1	B	302	ASP	2.9
1	F	263	ALA	2.9
1	E	112	ARG	2.9
1	F	205	ASP	2.9
1	D	61	ARG	2.8
1	F	28	GLN	2.8
1	B	112	ARG	2.8
1	B	26	ASP	2.8
1	B	70	ASP	2.8
1	D	94	GLN	2.8
1	C	7	LEU	2.7
1	D	26	ASP	2.7
1	E	157	TRP	2.6
1	C	112	ARG	2.6
1	D	7	LEU	2.6
1	E	205	ASP	2.5
1	F	26	ASP	2.5
1	B	7	LEU	2.5
1	A	3	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	E	7	LEU	2.4
1	C	168	GLU	2.4
1	E	28	GLN	2.4
1	C	157	TRP	2.4
1	D	21	VAL	2.4
1	C	61	ARG	2.4
1	E	94	GLN	2.4
1	F	22	GLN	2.4
1	D	105	ARG	2.4
1	F	72	LYS	2.4
1	D	76	GLN	2.3
1	E	196	ALA	2.3
1	A	26	ASP	2.3
1	C	197	ARG	2.3
1	B	168	GLU	2.3
1	E	3	GLY	2.3
1	B	20	LEU	2.3
1	D	132	ALA	2.3
1	A	76	GLN	2.3
1	E	214	VAL	2.3
1	A	7	LEU	2.3
1	D	266	THR	2.3
1	E	72	LYS	2.3
1	A	61	ARG	2.2
1	D	112	ARG	2.2
1	D	197	ARG	2.2
1	F	157	TRP	2.2
1	B	4	PRO	2.2
1	B	127	LEU	2.2
1	E	168	GLU	2.2
1	E	166	ALA	2.2
1	F	206	ASP	2.2
1	E	193	ALA	2.2
1	D	27	TRP	2.2
1	D	25	GLY	2.2
1	E	209	ALA	2.2
1	A	105	ARG	2.2
1	D	303	HIS	2.1
1	B	157	TRP	2.1
1	A	136	ASP	2.1
1	D	28	GLN	2.1
1	C	26	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	210	ASP	2.1
1	C	264	ASN	2.1
1	B	132	ALA	2.1
1	A	112	ARG	2.1
1	F	213	GLY	2.1
1	D	23	PRO	2.1
1	E	212	ALA	2.1
1	D	70	ASP	2.1
1	C	171	PRO	2.0
1	D	20	LEU	2.0
1	D	212	ALA	2.0
1	F	171	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	16D	B	1402	8/8	0.48	0.28	22,24,26,26	0
3	16D	E	1405	8/8	0.49	0.28	23,26,26,26	0
3	16D	C	1403	8/8	0.51	0.33	21,25,26,26	0
3	16D	F	1406	8/8	0.52	0.33	21,25,26,26	0
3	16D	D	1404	8/8	0.53	0.30	21,25,26,26	0
3	16D	A	1401	8/8	0.58	0.28	19,25,26,26	0
2	MN	E	1609	1/1	0.97	0.19	20,20,20,20	0
2	MN	B	1604	1/1	0.98	0.05	18,18,18,18	0
2	MN	E	1610	1/1	0.98	0.03	22,22,22,22	0
2	MN	F	1612	1/1	0.98	0.03	18,18,18,18	0
2	MN	C	1605	1/1	0.99	0.09	18,18,18,18	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MN	C	1606	1/1	0.99	0.07	17,17,17,17	0
2	MN	D	1607	1/1	0.99	0.10	20,20,20,20	0
2	MN	D	1608	1/1	0.99	0.02	18,18,18,18	0
2	MN	B	1603	1/1	0.99	0.09	18,18,18,18	0
2	MN	A	1602	1/1	0.99	0.02	16,16,16,16	0
2	MN	F	1611	1/1	0.99	0.11	16,16,16,16	0
2	MN	A	1601	1/1	1.00	0.12	12,12,12,12	0

6.5 Other polymers [i](#)

There are no such residues in this entry.