Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2021 – 07:18 PM EDT

PDB ID	:	1T5G
Title	:	Arginase-F2-L-Arginine complex
Authors	:	Cama, E.; Pethe, S.; Boucher, J.-L.; Han, S.; Emig, F.A.; Ash, D.E.; Viola, R.E.; Mansuy, D.; Christianson, D.W.
Deposited on	:	2004-05-04
Resolution	:	2.40 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 following versions of software and data (see references ⓘ) were used in the production of this report:

|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

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 2.40 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	Percentile Ranks	Value
Clashscore		21
Ramachandran outliers		1.0%
Sidechain outliers		6.2%

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution
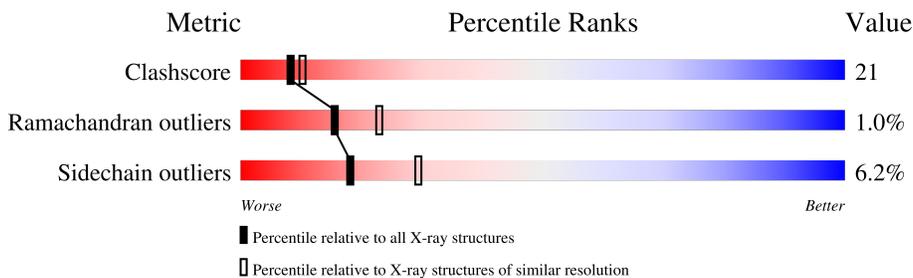
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	314	60% 36% •
1	B	314	60% 36% •
1	C	314	59% 38% •

WORLDWIDE
wwPDB
PROTEIN DATA BANK

2 Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 7300 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 Arginase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total 2384	C 1523	N 402	O 455	S 4	0	0	0
1	B	314	Total 2384	C 1523	N 402	O 455	S 4	0	0	0
1	C	314	Total 2384	C 1523	N 402	O 455	S 4	0	0	0

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	CYS	GLN	engineered mutation	UNP P07824
A	119	ALA	CYS	engineered mutation	UNP P07824
A	141	ALA	HIS	engineered mutation	UNP P07824
A	168	ALA	CYS	engineered mutation	UNP P07824
A	303	ALA	CYS	engineered mutation	UNP P07824
B	19	CYS	GLN	engineered mutation	UNP P07824
B	119	ALA	CYS	engineered mutation	UNP P07824
B	141	ALA	HIS	engineered mutation	UNP P07824
B	168	ALA	CYS	engineered mutation	UNP P07824
B	303	ALA	CYS	engineered mutation	UNP P07824
C	19	CYS	GLN	engineered mutation	UNP P07824
C	119	ALA	CYS	engineered mutation	UNP P07824
C	141	ALA	HIS	engineered mutation	UNP P07824
C	168	ALA	CYS	engineered mutation	UNP P07824
C	303	ALA	CYS	engineered mutation	UNP P07824

- Molecule 2 is FLUORIDE ION (three-letter code: F) (formula: F).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 2	F 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total F 2 2		0	0
2	C	2	Total F 2 2		0	0

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 4 is ARGININE (three-letter code: ARG) (formula: $C_6H_{15}N_4O_2$).
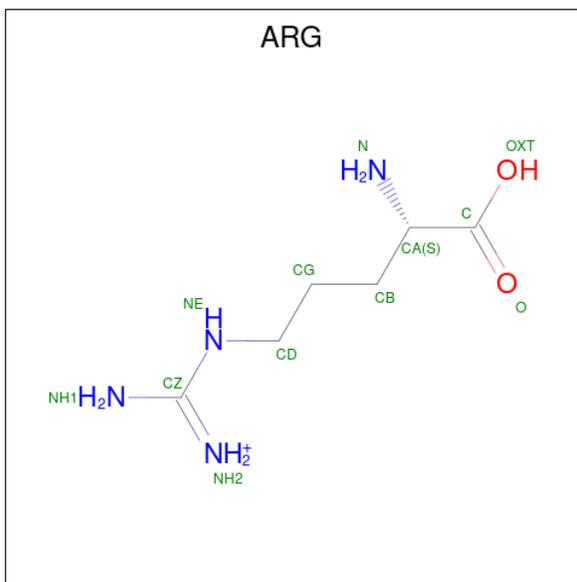
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total C N O 12 6 4 2				0	0
4	B	1	Total C N O 12 6 4 2				0	0
4	C	1	Total C N O 12 6 4 2				0	0

- Molecule 5 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	30	Total 30	O 30	0	0
5	B	32	Total 32	O 32	0	0
5	C	38	Total 38	O 38	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. 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. 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.

Note EDS was not executed.

• Molecule 1: Arginase 1

Chain A:

• Molecule 1: Arginase 1

Chain B:

• Molecule 1: Arginase 1

Chain C:
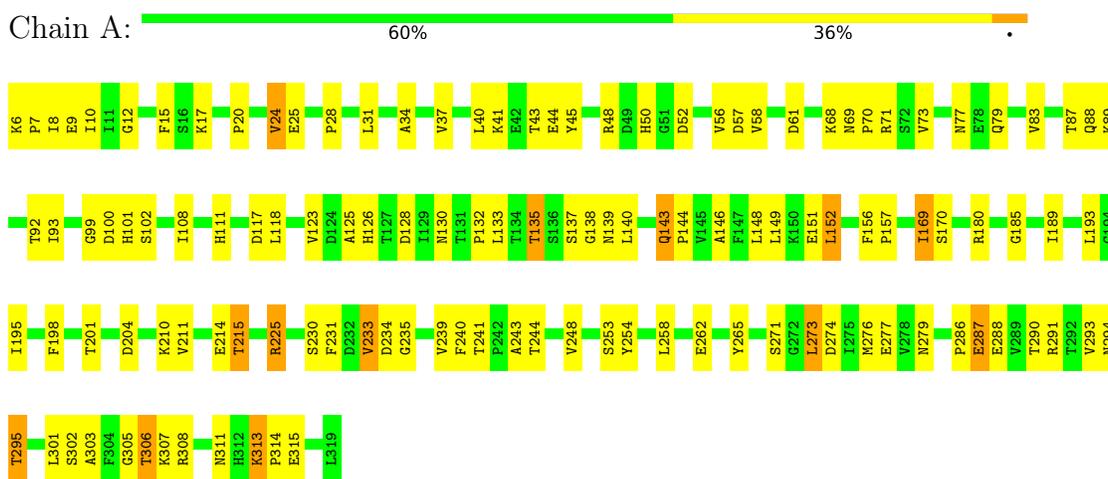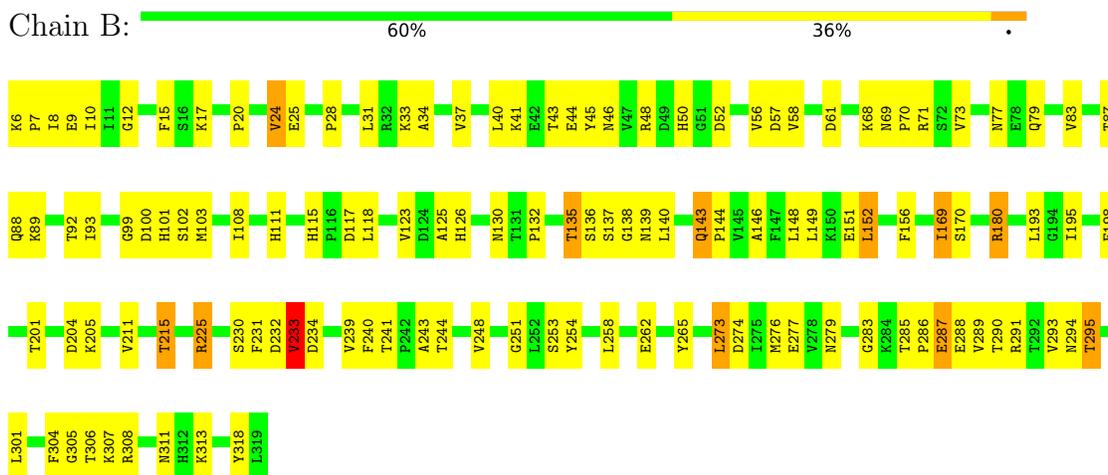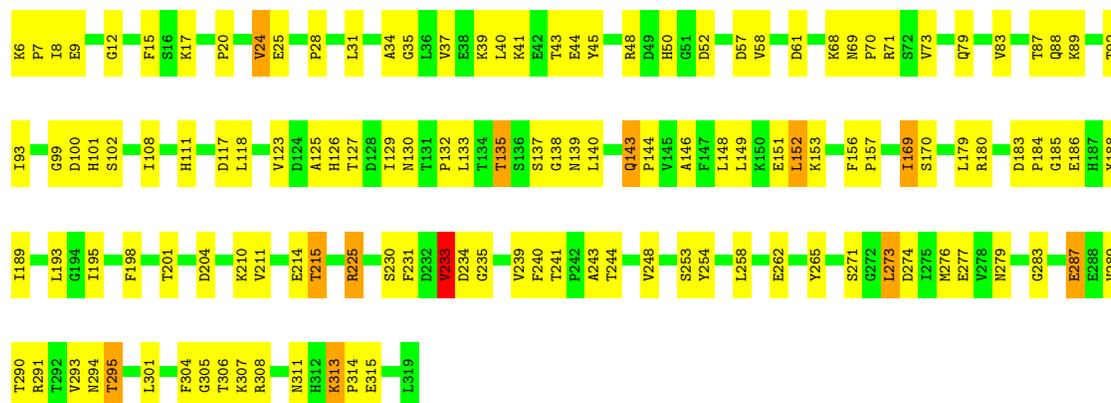
4 Data and refinement statistics ⓘ

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α, β, γ	87.72Å　87.72Å　104.52Å 90.00°　90.00°　120.00°	Depositor
Resolution (Å)	27.69　–　2.40	Depositor
% Data completeness (in resolution range)	95.0 (27.69-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.228　,　0.254	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7300	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

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: F, 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	0.40	0/2436	0.64	0/3309
1	B	0.40	0/2436	0.65	0/3309
1	C	0.41	0/2436	0.67	1/3309 (0.0%)
All	All	0.41	0/7308	0.65	1/9927 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	153	LYS	CD-CE-NZ	9.65	133.89	111.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	2384	0	2412	102	0
1	B	2384	0	2412	109	0
1	C	2384	0	2412	110	0
2	A	2	0	0	1	0
2	B	2	0	0	1	0
2	C	2	0	0	1	0
3	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
3	C	2	0	0	0	0
4	A	12	0	12	0	0
4	B	12	0	12	0	0
4	C	12	0	12	0	0
5	A	30	0	0	3	0
5	B	32	0	0	5	0
5	C	38	0	0	6	0
All	All	7300	0	7272	306	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 21.

All (306) 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:146:ALA:HA	1:C:152:LEU:HD23	1.55	0.89
1:A:146:ALA:HA	1:A:152:LEU:HD23	1.54	0.88
1:B:146:ALA:HA	1:B:152:LEU:HD23	1.55	0.87
1:C:135:THR:HB	1:C:137:SER:O	1.76	0.85
1:A:135:THR:HB	1:A:137:SER:O	1.77	0.85
1:B:135:THR:HB	1:B:137:SER:O	1.77	0.84
1:C:244:THR:HG22	1:C:277:GLU:O	1.76	0.84
1:C:126:HIS:HB3	5:C:743:HOH:O	1.79	0.83
1:B:9:GLU:OE1	1:B:48:ARG:HD3	1.81	0.81
1:A:15:PHE:CZ	1:A:17:LYS:HB2	2.17	0.80
1:A:244:THR:HG22	1:A:277:GLU:O	1.82	0.80
1:B:15:PHE:CZ	1:B:17:LYS:HB2	2.17	0.80
1:B:244:THR:HG22	1:B:277:GLU:O	1.81	0.80
1:C:15:PHE:CZ	1:C:17:LYS:HB2	2.17	0.80
1:A:9:GLU:OE1	1:A:48:ARG:HD3	1.82	0.80
1:B:151:GLU:OE2	1:B:170:SER:HB2	1.82	0.80
1:A:306:THR:HG21	5:A:722:HOH:O	1.83	0.79
1:B:143:GLN:H	1:B:144:PRO:HD3	1.47	0.79
1:C:9:GLU:OE1	1:C:48:ARG:HD3	1.82	0.78
1:C:151:GLU:OE2	1:C:170:SER:HB2	1.83	0.78
1:B:233:VAL:HG22	1:B:241:THR:HG21	1.64	0.77
1:A:151:GLU:OE2	1:A:170:SER:HB2	1.85	0.77
1:C:233:VAL:HG22	1:C:241:THR:HG21	1.67	0.76
1:C:307:LYS:H	1:C:311:ASN:HD21	1.34	0.76
1:A:126:HIS:HB3	5:A:702:HOH:O	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:SER:HA	1:B:144:PRO:HG3	1.69	0.74
1:A:307:LYS:H	1:A:311:ASN:HD21	1.36	0.74
1:C:102:SER:HA	1:C:144:PRO:HG3	1.70	0.74
1:B:307:LYS:H	1:B:311:ASN:HD21	1.36	0.73
1:A:79:GLN:O	1:A:83:VAL:HG23	1.88	0.73
1:B:125:ALA:HB2	1:B:234:ASP:O	1.86	0.73
1:A:102:SER:HA	1:A:144:PRO:HG3	1.71	0.72
1:A:40:LEU:HD23	1:A:301:LEU:HD23	1.71	0.72
1:C:40:LEU:HD23	1:C:301:LEU:HD23	1.71	0.72
1:C:198:PHE:CE2	1:C:215:THR:HG22	2.23	0.72
1:C:79:GLN:O	1:C:83:VAL:HG23	1.89	0.72
1:B:58:VAL:O	1:B:61:ASP:HB2	1.89	0.72
1:B:79:GLN:O	1:B:83:VAL:HG23	1.90	0.71
1:B:40:LEU:HD23	1:B:301:LEU:HD23	1.71	0.71
1:C:58:VAL:O	1:C:61:ASP:HB2	1.91	0.71
1:A:118:LEU:O	1:A:118:LEU:HD12	1.90	0.70
1:C:143:GLN:H	1:C:144:PRO:HD3	1.57	0.70
1:B:126:HIS:HB3	5:B:741:HOH:O	1.91	0.70
1:A:58:VAL:O	1:A:61:ASP:HB2	1.92	0.70
1:B:198:PHE:CE2	1:B:215:THR:HG22	2.26	0.69
1:A:198:PHE:CE2	1:A:215:THR:HG22	2.27	0.69
1:C:211:VAL:O	1:C:215:THR:HG23	1.91	0.69
1:B:211:VAL:O	1:B:215:THR:HG23	1.93	0.69
1:C:7:PRO:HB2	1:C:92:THR:HB	1.74	0.69
1:C:20:PRO:HD3	1:C:139:ASN:OD1	1.93	0.69
1:A:130:ASN:HB3	1:A:135:THR:HG23	1.74	0.69
1:A:211:VAL:O	1:A:215:THR:HG23	1.93	0.68
1:A:233:VAL:HG22	1:A:241:THR:HG21	1.76	0.68
1:B:143:GLN:N	1:B:144:PRO:CD	2.56	0.68
1:B:130:ASN:HB3	1:B:135:THR:HG23	1.76	0.68
1:B:7:PRO:HB2	1:B:92:THR:HB	1.76	0.67
1:C:100:ASP:HA	1:C:277:GLU:OE2	1.94	0.67
1:C:118:LEU:HD12	1:C:118:LEU:O	1.94	0.67
1:A:143:GLN:H	1:A:144:PRO:HD3	1.60	0.67
1:C:233:VAL:HG12	1:C:244:THR:HG21	1.77	0.67
1:A:233:VAL:HG12	1:A:244:THR:HG21	1.76	0.66
1:B:118:LEU:HD12	1:B:118:LEU:O	1.95	0.66
1:A:7:PRO:HB2	1:A:92:THR:HB	1.76	0.66
1:A:143:GLN:N	1:A:144:PRO:HD3	2.12	0.65
1:A:308:ARG:NH2	1:B:204:ASP:OD2	2.29	0.64
1:B:143:GLN:N	1:B:144:PRO:HD3	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:PRO:HD3	1:B:139:ASN:OD1	1.97	0.64
1:C:130:ASN:HB3	1:C:135:THR:HG23	1.79	0.64
1:C:143:GLN:N	1:C:144:PRO:CD	2.60	0.63
1:C:125:ALA:HB2	1:C:234:ASP:O	2.00	0.62
1:A:143:GLN:N	1:A:144:PRO:CD	2.63	0.62
1:C:143:GLN:N	1:C:144:PRO:HD3	2.14	0.62
1:B:233:VAL:HG12	1:B:244:THR:HG21	1.80	0.61
1:C:24:VAL:HB	1:C:99:GLY:CA	2.30	0.61
1:B:225:ARG:HG3	1:B:225:ARG:HH11	1.65	0.61
1:A:8:ILE:HD13	1:A:93:ILE:HB	1.82	0.61
1:A:20:PRO:HD3	1:A:139:ASN:OD1	2.01	0.61
1:B:152:LEU:HD13	1:B:193:LEU:HD21	1.83	0.60
1:B:8:ILE:HD13	1:B:93:ILE:HB	1.83	0.60
1:B:273:LEU:HD23	1:B:274:ASP:N	2.16	0.60
1:A:24:VAL:HB	1:A:99:GLY:CA	2.30	0.60
1:B:100:ASP:HA	1:B:277:GLU:OE2	2.01	0.60
1:A:254:TYR:CE1	1:A:295:THR:HB	2.38	0.59
1:A:12:GLY:HA3	1:A:52:ASP:OD1	2.03	0.59
1:C:152:LEU:HD13	1:C:193:LEU:HD21	1.83	0.59
1:A:273:LEU:HD23	1:A:274:ASP:N	2.17	0.59
1:B:24:VAL:HB	1:B:99:GLY:CA	2.32	0.59
1:C:254:TYR:CE1	1:C:295:THR:HB	2.38	0.58
1:B:225:ARG:HG3	1:B:225:ARG:NH1	2.18	0.58
1:C:273:LEU:HD23	1:C:274:ASP:N	2.18	0.58
1:A:225:ARG:HG3	1:A:225:ARG:HH11	1.69	0.58
1:C:73:VAL:HG21	1:C:140:LEU:HD12	1.86	0.58
1:A:100:ASP:HA	1:A:277:GLU:OE2	2.03	0.58
1:C:243:ALA:HB1	1:C:279:ASN:O	2.04	0.58
1:A:243:ALA:HB1	1:A:279:ASN:O	2.03	0.57
1:B:117:ASP:O	1:B:225:ARG:HG3	2.04	0.57
1:B:232:ASP:O	1:B:234:ASP:N	2.38	0.57
1:B:232:ASP:C	1:B:234:ASP:H	2.07	0.57
1:B:205:LYS:HG3	5:B:707:HOH:O	2.05	0.57
1:A:254:TYR:HE1	1:A:295:THR:HB	1.69	0.56
1:B:140:LEU:O	1:B:144:PRO:HD3	2.05	0.56
1:C:8:ILE:HD13	1:C:93:ILE:HB	1.86	0.56
1:A:73:VAL:HG21	1:A:140:LEU:HD12	1.87	0.56
1:B:254:TYR:CE1	1:B:295:THR:HB	2.41	0.56
1:A:152:LEU:HD13	1:A:193:LEU:HD21	1.86	0.56
1:C:254:TYR:HE1	1:C:295:THR:HB	1.70	0.56
1:B:132:PRO:HD2	1:B:156:PHE:CD2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:273:LEU:HD23	1:B:274:ASP:H	1.69	0.56
1:C:225:ARG:HG3	1:C:225:ARG:HH11	1.71	0.55
1:A:152:LEU:CD1	1:A:193:LEU:HD21	2.36	0.55
1:B:12:GLY:HA3	1:B:52:ASP:OD1	2.05	0.55
1:B:169:ILE:HD13	1:B:169:ILE:H	1.72	0.55
1:C:273:LEU:HD23	1:C:274:ASP:H	1.71	0.55
1:A:225:ARG:HG3	1:A:225:ARG:NH1	2.21	0.55
1:C:12:GLY:HA3	1:C:52:ASP:OD1	2.06	0.55
1:A:273:LEU:HD23	1:A:274:ASP:H	1.71	0.55
1:C:258:LEU:O	1:C:262:GLU:HG3	2.07	0.55
1:A:306:THR:CG2	5:A:722:HOH:O	2.50	0.55
1:A:43:THR:C	1:A:45:TYR:H	2.10	0.55
1:A:24:VAL:HB	1:A:99:GLY:HA2	1.90	0.54
1:A:169:ILE:HD13	1:A:169:ILE:H	1.71	0.54
1:A:308:ARG:HB2	1:B:201:THR:HG23	1.90	0.54
1:B:311:ASN:O	1:C:184:PRO:HA	2.08	0.54
1:C:225:ARG:HG3	1:C:225:ARG:NH1	2.23	0.54
1:C:132:PRO:HD2	1:C:156:PHE:CD2	2.42	0.54
1:C:140:LEU:O	1:C:144:PRO:HD3	2.07	0.54
1:C:117:ASP:O	1:C:225:ARG:HG3	2.08	0.53
1:B:101:HIS:CD2	1:B:276:MET:HG3	2.43	0.53
1:C:186:GLU:HG3	5:C:729:HOH:O	2.07	0.53
1:A:117:ASP:O	1:A:225:ARG:HG3	2.08	0.53
1:B:43:THR:C	1:B:45:TYR:H	2.11	0.53
1:B:243:ALA:HB1	1:B:279:ASN:O	2.09	0.53
1:C:43:THR:C	1:C:45:TYR:H	2.10	0.53
1:C:152:LEU:CD1	1:C:193:LEU:HD21	2.39	0.53
1:B:254:TYR:HE1	1:B:295:THR:HB	1.72	0.53
1:C:169:ILE:HD13	1:C:169:ILE:H	1.74	0.53
1:B:68:LYS:HE3	1:B:138:GLY:H	1.73	0.53
1:B:73:VAL:HG21	1:B:140:LEU:HD12	1.90	0.53
1:C:24:VAL:HB	1:C:99:GLY:HA2	1.90	0.53
1:B:152:LEU:CD1	1:B:193:LEU:HD21	2.38	0.53
1:B:24:VAL:HB	1:B:99:GLY:HA2	1.92	0.52
1:C:233:VAL:CG1	1:C:244:THR:HG21	2.39	0.52
1:B:33:LYS:HE2	5:B:754:HOH:O	2.09	0.52
1:C:287:GLU:HG2	5:C:787:HOH:O	2.09	0.52
1:C:31:LEU:HD23	1:C:293:VAL:HG13	1.91	0.52
1:B:31:LEU:HD23	1:B:293:VAL:HG13	1.91	0.52
1:C:68:LYS:HE3	1:C:138:GLY:H	1.73	0.52
1:A:50:HIS:O	1:A:50:HIS:ND1	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:LYS:HE3	1:A:138:GLY:H	1.75	0.52
1:A:233:VAL:HG12	1:A:244:THR:CG2	2.40	0.52
1:A:132:PRO:HD2	1:A:156:PHE:CD2	2.45	0.51
1:C:101:HIS:CD2	1:C:276:MET:HG3	2.45	0.51
1:B:308:ARG:NH2	1:C:204:ASP:OD2	2.41	0.51
1:C:9:GLU:OE2	1:C:50:HIS:HB2	2.10	0.50
1:A:307:LYS:H	1:A:311:ASN:ND2	2.08	0.50
1:A:101:HIS:CD2	1:A:276:MET:HG3	2.47	0.50
1:B:233:VAL:HG13	1:B:233:VAL:O	2.12	0.50
1:B:318:TYR:CD1	1:C:188:TYR:CG	3.00	0.50
1:B:25:GLU:O	1:B:28:PRO:HD2	2.12	0.50
1:A:70:PRO:HG2	1:A:71:ARG:H	1.77	0.49
1:B:50:HIS:O	1:B:50:HIS:ND1	2.45	0.49
1:C:6:LYS:HD2	1:C:7:PRO:HD3	1.94	0.49
1:A:6:LYS:HD2	1:A:7:PRO:HD3	1.94	0.49
1:A:125:ALA:HB2	1:A:234:ASP:O	2.12	0.49
1:B:240:PHE:CE2	1:B:253:SER:HA	2.48	0.49
1:A:240:PHE:CE2	1:A:253:SER:HA	2.47	0.49
1:B:70:PRO:HG2	1:B:71:ARG:H	1.77	0.49
1:B:307:LYS:H	1:B:311:ASN:ND2	2.08	0.49
1:B:308:ARG:HE	1:C:201:THR:HG22	1.76	0.49
1:C:307:LYS:H	1:C:311:ASN:ND2	2.06	0.49
1:A:31:LEU:HD23	1:A:293:VAL:HG13	1.95	0.48
1:A:69:ASN:O	1:A:73:VAL:HG23	2.13	0.48
1:A:128:ASP:HB3	1:A:144:PRO:HD2	1.96	0.48
1:B:89:LYS:HB2	1:B:89:LYS:NZ	2.29	0.48
1:B:262:GLU:OE1	1:C:201:THR:HG22	2.12	0.48
1:C:240:PHE:CE2	1:C:253:SER:HA	2.48	0.48
1:B:232:ASP:C	1:B:234:ASP:N	2.65	0.48
1:C:233:VAL:HG12	1:C:244:THR:CG2	2.44	0.48
1:B:233:VAL:HG12	1:B:244:THR:CG2	2.44	0.48
1:C:265:TYR:CZ	1:C:305:GLY:HA2	2.47	0.48
1:B:61:ASP:OD1	1:B:71:ARG:HB2	2.14	0.47
1:B:308:ARG:HB2	1:C:201:THR:HG23	1.95	0.47
1:C:239:VAL:HG23	5:C:712:HOH:O	2.13	0.47
1:C:180:ARG:NH2	1:C:235:GLY:O	2.47	0.47
1:A:273:LEU:CD2	1:A:274:ASP:N	2.77	0.47
1:C:273:LEU:CD2	1:C:274:ASP:N	2.77	0.47
1:A:265:TYR:CZ	1:A:305:GLY:HA2	2.50	0.47
1:B:273:LEU:CD2	1:B:274:ASP:N	2.77	0.47
1:C:230:SER:HA	1:C:274:ASP:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:LYS:HB2	1:A:89:LYS:NZ	2.30	0.46
1:A:118:LEU:HD12	1:A:118:LEU:C	2.36	0.46
1:C:89:LYS:HB2	1:C:89:LYS:NZ	2.30	0.46
1:C:101:HIS:N	1:C:277:GLU:OE2	2.47	0.46
1:C:70:PRO:HG2	1:C:71:ARG:H	1.78	0.46
1:B:37:VAL:O	1:B:41:LYS:HG3	2.16	0.46
1:C:20:PRO:HD3	1:C:139:ASN:CG	2.36	0.46
1:C:69:ASN:O	1:C:73:VAL:HG23	2.16	0.46
1:B:233:VAL:CG1	1:B:244:THR:HG21	2.45	0.46
1:B:108:ILE:HD12	1:B:148:LEU:CD1	2.46	0.46
1:A:189:ILE:O	1:A:193:LEU:HD13	2.16	0.46
1:C:108:ILE:HD12	1:C:148:LEU:CD1	2.46	0.45
1:B:262:GLU:OE1	1:C:201:THR:CG2	2.65	0.45
1:A:20:PRO:HD3	1:A:139:ASN:CG	2.37	0.45
1:A:290:THR:HG22	1:A:294:ASN:ND2	2.32	0.45
1:B:136:SER:HB2	5:B:799:HOH:O	2.16	0.45
1:C:101:HIS:HB2	2:C:601:F:F	2.07	0.45
1:B:230:SER:HA	1:B:274:ASP:HB2	1.99	0.45
1:C:290:THR:HG22	1:C:294:ASN:ND2	2.32	0.45
1:A:88:GLN:HG3	1:A:111:HIS:HA	1.98	0.45
1:A:239:VAL:HG12	1:A:239:VAL:O	2.15	0.45
1:A:108:ILE:HD12	1:A:148:LEU:CD1	2.47	0.44
1:C:287:GLU:CG	5:C:787:HOH:O	2.64	0.44
1:A:146:ALA:HA	1:A:152:LEU:CD2	2.36	0.44
1:A:169:ILE:HD13	1:A:169:ILE:N	2.32	0.44
1:B:20:PRO:HD3	1:B:139:ASN:CG	2.37	0.44
1:A:140:LEU:O	1:A:144:PRO:HD3	2.18	0.44
1:B:265:TYR:CZ	1:B:305:GLY:HA2	2.52	0.44
1:B:291:ARG:O	1:B:295:THR:CG2	2.65	0.44
1:C:118:LEU:HD12	1:C:118:LEU:C	2.38	0.44
1:B:69:ASN:O	1:B:73:VAL:HG23	2.18	0.44
1:B:101:HIS:HB2	2:B:605:F:F	2.08	0.44
1:B:149:LEU:HD11	1:B:195:ILE:HD11	1.99	0.44
1:C:61:ASP:OD1	1:C:71:ARG:HB2	2.18	0.44
1:A:123:VAL:HG11	1:A:231:PHE:CZ	2.53	0.44
1:A:201:THR:HG23	1:C:308:ARG:HB2	1.98	0.44
1:A:308:ARG:O	1:B:180:ARG:HA	2.17	0.44
1:A:308:ARG:HB2	1:B:201:THR:CG2	2.48	0.44
1:A:291:ARG:O	1:A:295:THR:CG2	2.66	0.44
1:C:88:GLN:HG3	1:C:111:HIS:HA	2.00	0.44
1:C:233:VAL:CG1	1:C:244:THR:CG2	2.96	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ARG:NH2	1:A:235:GLY:O	2.51	0.43
1:B:130:ASN:O	1:B:143:GLN:HG2	2.18	0.43
1:C:25:GLU:O	1:C:28:PRO:HD2	2.19	0.43
1:A:230:SER:HA	1:A:274:ASP:HB2	2.00	0.43
1:C:149:LEU:HD11	1:C:195:ILE:HD11	2.00	0.43
1:A:180:ARG:HH11	1:A:248:VAL:HB	1.84	0.43
1:A:204:ASP:OD2	1:C:308:ARG:NH2	2.48	0.43
1:A:25:GLU:O	1:A:28:PRO:HD2	2.19	0.43
1:A:61:ASP:OD1	1:A:71:ARG:HB2	2.17	0.43
1:B:123:VAL:HG11	1:B:231:PHE:CZ	2.54	0.43
1:B:287:GLU:HG3	1:B:288:GLU:N	2.34	0.43
1:A:287:GLU:HG3	1:A:288:GLU:N	2.34	0.43
1:B:88:GLN:HG3	1:B:111:HIS:HA	2.01	0.43
1:C:189:ILE:O	1:C:193:LEU:HD13	2.18	0.43
1:B:56:VAL:HG23	1:B:56:VAL:O	2.19	0.43
1:A:101:HIS:HB2	2:A:603:F:F	2.09	0.43
1:A:185:GLY:O	1:A:189:ILE:HG13	2.19	0.43
1:C:34:ALA:HB2	1:C:293:VAL:HG11	2.01	0.43
1:C:180:ARG:HH11	1:C:248:VAL:HB	1.83	0.43
1:B:6:LYS:HA	1:B:7:PRO:HD3	1.79	0.42
1:B:8:ILE:HD12	1:B:304:PHE:CE1	2.54	0.42
1:C:313:LYS:HA	1:C:314:PRO:HD3	1.84	0.42
1:A:10:ILE:HD12	1:A:10:ILE:N	2.34	0.42
1:A:117:ASP:O	1:A:225:ARG:NH1	2.53	0.42
1:B:239:VAL:HG12	1:B:239:VAL:O	2.19	0.42
1:C:169:ILE:HD13	1:C:169:ILE:N	2.33	0.42
1:B:6:LYS:HD2	1:B:7:PRO:HD3	2.01	0.42
1:B:262:GLU:OE2	1:B:308:ARG:HG3	2.19	0.42
1:C:35:GLY:O	1:C:39:LYS:HG3	2.19	0.42
1:A:93:ILE:HG23	1:A:271:SER:O	2.19	0.42
1:C:37:VAL:O	1:C:41:LYS:HG3	2.20	0.42
1:C:239:VAL:O	1:C:239:VAL:HG12	2.18	0.42
1:C:265:TYR:CE1	1:C:305:GLY:HA2	2.55	0.42
1:B:180:ARG:HH11	1:B:248:VAL:HB	1.85	0.42
1:B:283:GLY:HA3	1:B:289:VAL:HG23	2.02	0.42
1:A:258:LEU:O	1:A:262:GLU:HG3	2.19	0.42
1:B:37:VAL:HG12	1:B:41:LYS:HE3	2.00	0.42
1:C:183:ASP:HB2	5:C:729:HOH:O	2.19	0.42
1:A:233:VAL:CG1	1:A:244:THR:HG21	2.46	0.42
1:B:290:THR:HG22	1:B:294:ASN:ND2	2.35	0.42
1:C:6:LYS:HA	1:C:7:PRO:HD3	1.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:VAL:O	1:A:41:LYS:HG3	2.18	0.42
1:B:10:ILE:N	1:B:10:ILE:HD12	2.35	0.42
1:C:291:ARG:O	1:C:295:THR:CG2	2.68	0.42
1:A:6:LYS:HG2	1:A:93:ILE:HD11	2.01	0.41
1:A:314:PRO:O	1:A:315:GLU:HB2	2.19	0.41
1:C:37:VAL:HG12	1:C:41:LYS:HE3	2.02	0.41
1:A:34:ALA:HB2	1:A:293:VAL:HG11	2.01	0.41
1:A:73:VAL:HG12	1:A:77:ASN:ND2	2.35	0.41
1:A:56:VAL:HG23	1:A:56:VAL:O	2.20	0.41
1:A:133:LEU:HG	1:A:157:PRO:HD2	2.03	0.41
1:B:68:LYS:HG3	1:B:138:GLY:HA3	2.01	0.41
1:B:118:LEU:HD12	1:B:118:LEU:C	2.39	0.41
1:B:34:ALA:HB2	1:B:293:VAL:HG11	2.02	0.41
1:A:210:LYS:HE3	1:A:214:GLU:HG3	2.01	0.41
1:B:308:ARG:O	1:C:180:ARG:HA	2.21	0.41
1:C:210:LYS:HE3	1:C:214:GLU:HG3	2.02	0.41
1:B:318:TYR:CD2	1:C:185:GLY:HA2	2.56	0.41
1:A:37:VAL:HG12	1:A:41:LYS:HE3	2.02	0.41
1:A:149:LEU:HD11	1:A:195:ILE:HD11	2.01	0.41
1:A:302:SER:O	1:A:303:ALA:C	2.59	0.41
1:B:258:LEU:O	1:B:262:GLU:HG3	2.21	0.41
1:C:123:VAL:HG11	1:C:231:PHE:CZ	2.56	0.41
1:C:129:ILE:HG12	1:C:129:ILE:O	2.21	0.41
1:C:234:ASP:OD1	1:C:234:ASP:C	2.58	0.41
1:A:313:LYS:HA	1:A:314:PRO:HD3	1.83	0.41
1:B:115:HIS:CE1	5:B:717:HOH:O	2.74	0.41
1:C:127:THR:HG23	1:C:179:LEU:HD22	2.01	0.40
1:B:77:ASN:OD1	1:B:103:MET:HA	2.22	0.40
1:C:93:ILE:HG23	1:C:271:SER:O	2.21	0.40
1:C:283:GLY:HA3	1:C:289:VAL:HG23	2.02	0.40
1:C:314:PRO:O	1:C:315:GLU:HB2	2.21	0.40
1:C:8:ILE:HD12	1:C:304:PHE:CE1	2.56	0.40
1:B:180:ARG:NE	1:B:251:GLY:HA2	2.37	0.40
1:B:285:THR:OG1	1:B:288:GLU:HG3	2.21	0.40
1:C:133:LEU:HG	1:C:157:PRO:HD2	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	312/314 (99%)	289 (93%)	21 (7%)	2 (1%)	25	36
1	B	312/314 (99%)	292 (94%)	16 (5%)	4 (1%)	12	17
1	C	312/314 (99%)	291 (93%)	18 (6%)	3 (1%)	15	23
All	All	936/942 (99%)	872 (93%)	55 (6%)	9 (1%)	15	23

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	233	VAL
1	A	44	GLU
1	B	44	GLU
1	B	143	GLN
1	C	44	GLU
1	A	143	GLN
1	B	180	ARG
1	C	143	GLN
1	C	233	VAL

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	259/260 (100%)	244 (94%)	15 (6%)	20	32
1	B	259/260 (100%)	243 (94%)	16 (6%)	18	29
1	C	259/260 (100%)	245 (95%)	14 (5%)	22	36

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	777/780 (100%)	732 (94%)	45 (6%)	18 32

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	VAL
1	A	57	ASP
1	A	87	THR
1	A	135	THR
1	A	152	LEU
1	A	169	ILE
1	A	215	THR
1	A	225	ARG
1	A	233	VAL
1	A	273	LEU
1	A	286	PRO
1	A	287	GLU
1	A	295	THR
1	A	306	THR
1	A	313	LYS
1	B	24	VAL
1	B	46	ASN
1	B	57	ASP
1	B	87	THR
1	B	135	THR
1	B	152	LEU
1	B	169	ILE
1	B	215	THR
1	B	225	ARG
1	B	233	VAL
1	B	273	LEU
1	B	286	PRO
1	B	287	GLU
1	B	295	THR
1	B	306	THR
1	B	313	LYS
1	C	24	VAL
1	C	57	ASP
1	C	87	THR
1	C	135	THR
1	C	152	LEU
1	C	169	ILE

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Mol	Chain	Res	Type
1	C	215	THR
1	C	225	ARG
1	C	233	VAL
1	C	273	LEU
1	C	287	GLU
1	C	295	THR
1	C	306	THR
1	C	313	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	88	GLN
1	A	115	HIS
1	A	311	ASN
1	B	79	GLN
1	B	88	GLN
1	B	115	HIS
1	B	311	ASN
1	C	79	GLN
1	C	88	GLN
1	C	115	HIS
1	C	311	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 15 ligands modelled in this entry, 12 are monoatomic - leaving 3 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	#$|Z| > 2$
4	ARG	C	1002	-	7,11,11	1.16	1 (14%)	6,13,13	0.60	0
4	ARG	B	1001	-	7,11,11	1.10	1 (14%)	6,13,13	0.75	0
4	ARG	A	1000	-	7,11,11	1.16	1 (14%)	6,13,13	0.79	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
4	ARG	C	1002	-	-	3/7/11/11	-
4	ARG	B	1001	-	-	4/7/11/11	-
4	ARG	A	1000	-	-	4/7/11/11	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1000	ARG	CD-NE	2.27	1.51	1.46
4	B	1001	ARG	CD-NE	2.06	1.50	1.46
4	C	1002	ARG	CD-NE	2.01	1.50	1.46

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1000	ARG	C-CA-CB-CG
4	B	1001	ARG	C-CA-CB-CG

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Mol	Chain	Res	Type	Atoms
4	C	1002	ARG	C-CA-CB-CG
4	C	1002	ARG	NE-CD-CG-CB
4	B	1001	ARG	NE-CD-CG-CB
4	A	1000	ARG	NE-CD-CG-CB
4	A	1000	ARG	N-CA-CB-CG
4	B	1001	ARG	N-CA-CB-CG
4	C	1002	ARG	N-CA-CB-CG
4	B	1001	ARG	CA-CB-CG-CD
4	A	1000	ARG	CA-CB-CG-CD

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.