



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 03:58 am BST

PDB ID : 3UT3
Title : A novel PAI-I inhibitor and its structural mechanism
Authors : Lin, Z.H.; Hong, Z.B.; Shi, X.L.; Hu, L.H.; Andreasen, P.A.; Huang, M.D.
Deposited on : 2011-11-25
Resolution : 2.42 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) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

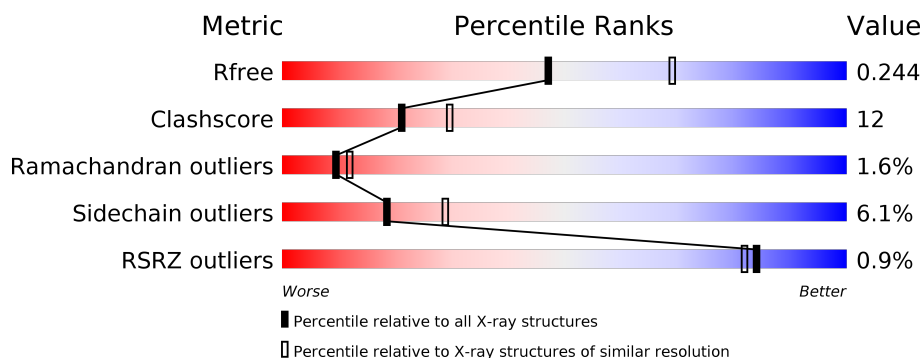
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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}	130704	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

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 on 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	375	<div> <div style="width: 67%;"></div> <div style="width: 27%;"></div> <div style="width: 6%;"></div> <div style="width: 2%;"></div> <div style="width: 2%;"></div> </div>
1	B	375	<div> <div style="width: 67%;"></div> <div style="width: 26%;"></div> <div style="width: 6%;"></div> <div style="width: 2%;"></div> <div style="width: 2%;"></div> </div>
1	C	375	<div> <div style="width: 74%;"></div> <div style="width: 22%;"></div> <div style="width: 6%;"></div> <div style="width: 2%;"></div> <div style="width: 2%;"></div> </div>
1	D	375	<div> <div style="width: 69%;"></div> <div style="width: 23%;"></div> <div style="width: 6%;"></div> <div style="width: 2%;"></div> <div style="width: 2%;"></div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11638 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.

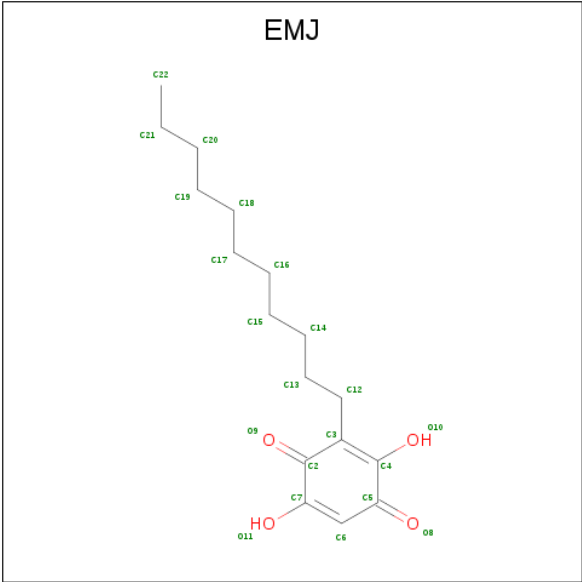
- Molecule 1 is a protein called Plasminogen activator inhibitor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	0	0	0
			2928	1881	499	533	15			
1	B	360	Total	C	N	O	S	0	0	0
			2877	1849	489	525	14			
1	C	368	Total	C	N	O	S	0	0	0
			2935	1886	500	534	15			
1	D	360	Total	C	N	O	S	0	0	0
			2877	1849	489	525	14			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	150	HIS	ASN	ENGINEERED MUTATION	UNP P05121
A	154	THR	LYS	ENGINEERED MUTATION	UNP P05121
A	319	LEU	GLN	ENGINEERED MUTATION	UNP P05121
A	354	ILE	MET	ENGINEERED MUTATION	UNP P05121
B	150	HIS	ASN	ENGINEERED MUTATION	UNP P05121
B	154	THR	LYS	ENGINEERED MUTATION	UNP P05121
B	319	LEU	GLN	ENGINEERED MUTATION	UNP P05121
B	354	ILE	MET	ENGINEERED MUTATION	UNP P05121
C	150	HIS	ASN	ENGINEERED MUTATION	UNP P05121
C	154	THR	LYS	ENGINEERED MUTATION	UNP P05121
C	319	LEU	GLN	ENGINEERED MUTATION	UNP P05121
C	354	ILE	MET	ENGINEERED MUTATION	UNP P05121
D	150	HIS	ASN	ENGINEERED MUTATION	UNP P05121
D	154	THR	LYS	ENGINEERED MUTATION	UNP P05121
D	319	LEU	GLN	ENGINEERED MUTATION	UNP P05121
D	354	ILE	MET	ENGINEERED MUTATION	UNP P05121

- Molecule 2 is 2,5-dihydroxy-3-undecylcyclohexa-2,5-diene-1,4-dione (three-letter code: EMJ) (formula: C₁₇H₂₆O₄).

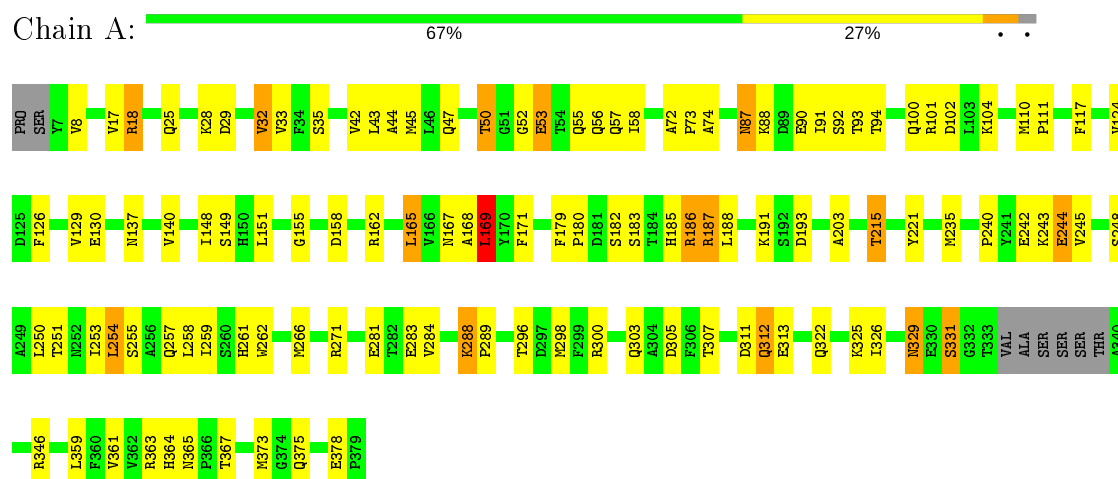


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			21	17	4		

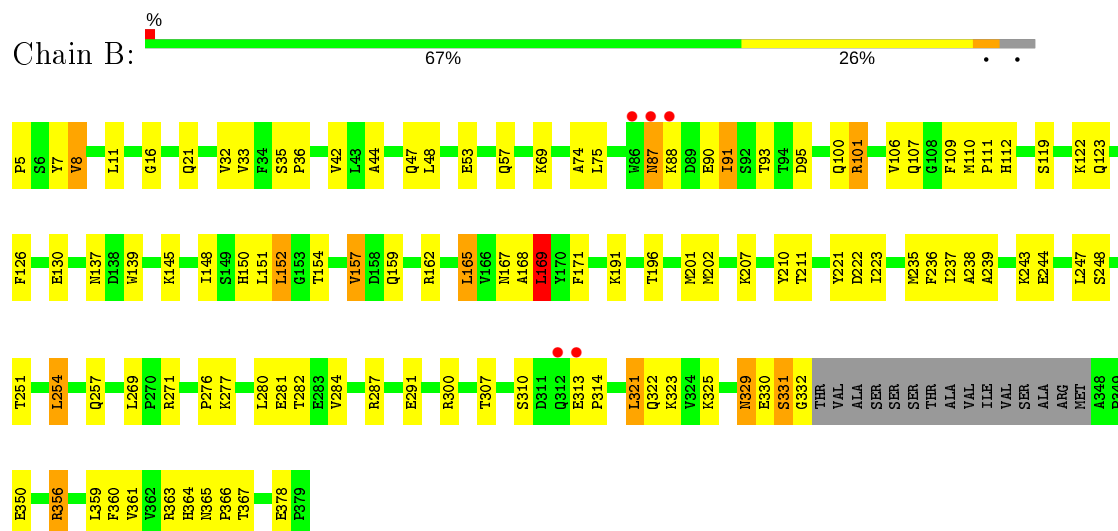
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

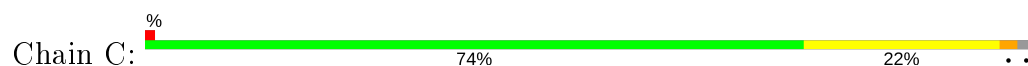
• Molecule 1: Plasminogen activator inhibitor 1

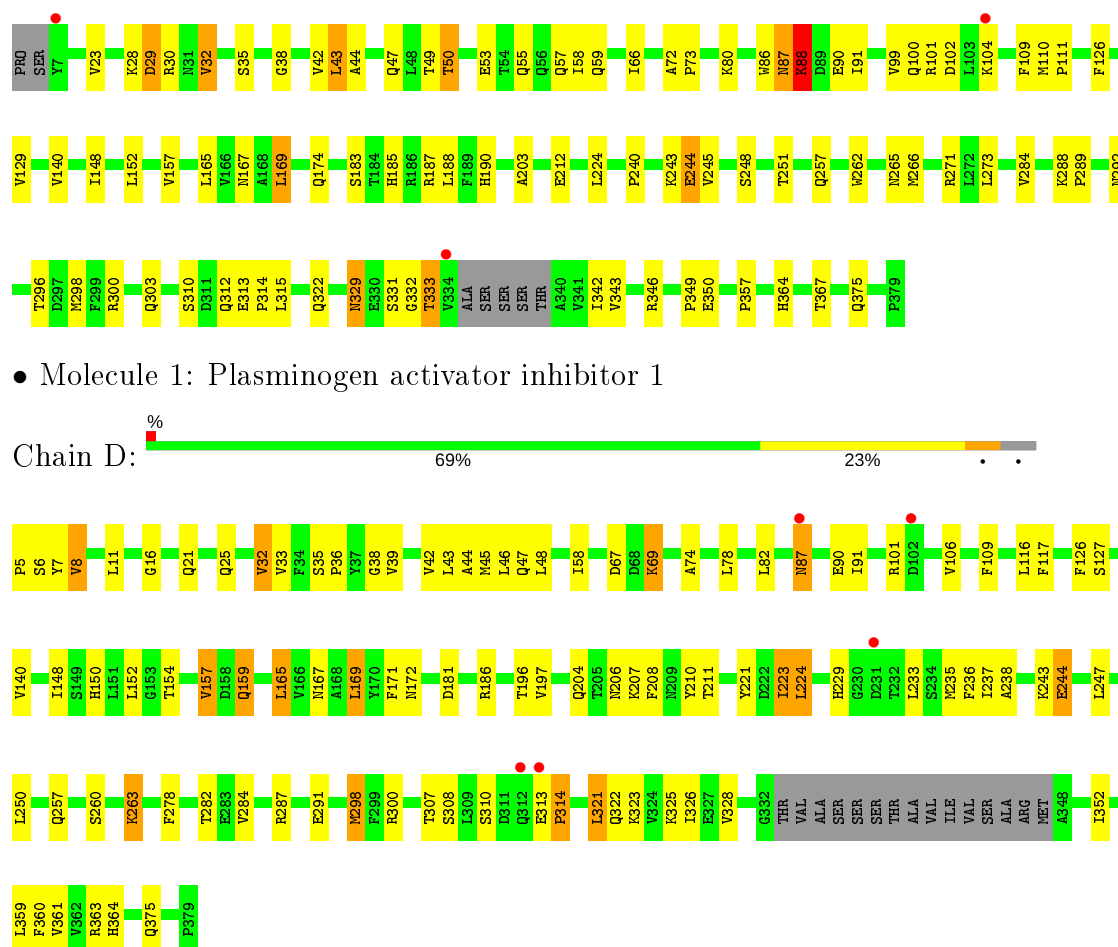


• Molecule 1: Plasminogen activator inhibitor 1



• Molecule 1: Plasminogen activator inhibitor 1





• Molecule 1: Plasminogen activator inhibitor 1

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	65.25Å 74.95Å 103.78Å 90.89° 93.33° 115.72°	Depositor
Resolution (Å)	58.69 – 2.42 57.72 – 2.42	Depositor EDS
% Data completeness (in resolution range)	96.6 (58.69-2.42) 96.6 (57.72-2.42)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.202 , 0.245 0.206 , 0.244	Depositor DCC
R_{free} test set	3288 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	34.5	Xtriage
Anisotropy	0.119	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.008 for -h,h+k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11638	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.22% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.75	1/2998 (0.0%)	0.86	5/4063 (0.1%)
1	B	0.73	0/2948	0.85	3/3995 (0.1%)
1	C	0.74	0/3005	0.83	2/4073 (0.0%)
1	D	0.72	0/2948	0.83	2/3995 (0.1%)
All	All	0.74	1/11899 (0.0%)	0.84	12/16126 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	283	GLU	CD-OE1	5.01	1.31	1.25

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	169	LEU	CA-CB-CG	7.55	132.66	115.30
1	A	169	LEU	CA-CB-CG	7.11	131.65	115.30
1	A	18	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	A	187	ARG	NE-CZ-NH1	6.18	123.39	120.30
1	D	321	LEU	CA-CB-CG	6.03	129.17	115.30
1	D	169	LEU	CA-CB-CG	5.81	128.67	115.30
1	A	235	MET	CG-SD-CE	-5.77	90.97	100.20
1	B	321	LEU	CA-CB-CG	5.63	128.25	115.30
1	C	43	LEU	CA-CB-CG	-5.59	102.45	115.30
1	B	169	LEU	CA-CB-CG	5.37	127.66	115.30
1	A	288	LYS	CD-CE-NZ	5.18	123.61	111.70
1	B	162	ARG	NE-CZ-NH1	5.07	122.84	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2928	0	2925	75	0
1	B	2877	0	2865	80	0
1	C	2935	0	2934	65	0
1	D	2877	0	2865	67	0
2	B	21	0	25	7	0
All	All	11638	0	11614	281	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 12.

All (281) 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:42:VAL:HG21	1:C:167:ASN:HB2	1.52	0.90
1:C:35:SER:H	1:C:322:GLN:HE22	1.21	0.89
1:A:35:SER:H	1:A:322:GLN:HE22	1.19	0.88
1:A:191:LYS:HD3	1:A:378:GLU:OE1	1.73	0.88
1:A:42:VAL:HG21	1:A:167:ASN:HB2	1.59	0.83
1:C:87:ASN:O	1:C:88:LYS:HB3	1.79	0.80
1:C:174:GLN:HG2	1:C:332:GLY:HA3	1.65	0.79
1:D:243:LYS:O	1:D:244:GLU:HB2	1.85	0.76
1:B:221:TYR:CE1	1:B:223:ILE:HD11	2.22	0.74
1:B:247:LEU:HD11	1:B:359:LEU:HD12	1.68	0.74
1:A:17:VAL:HG22	1:A:373:MET:CE	2.17	0.74
1:C:329:ASN:ND2	1:C:331:SER:O	2.24	0.71
1:C:87:ASN:O	1:C:88:LYS:CB	2.37	0.70
1:C:243:LYS:O	1:C:244:GLU:HB2	1.92	0.70
1:A:58:ILE:HD11	1:A:298:MET:HE2	1.74	0.69
1:B:364:HIS:HD2	1:B:367:THR:OG1	1.76	0.69
1:C:87:ASN:O	1:C:90:GLU:OE2	2.14	0.65
1:A:17:VAL:HG22	1:A:373:MET:HE3	1.79	0.65
1:D:78:LEU:O	1:D:82:LEU:HG	1.97	0.65
1:B:251:THR:HG23	1:B:359:LEU:HD11	1.78	0.64
1:C:50:THR:HG23	1:C:55:GLN:HB2	1.81	0.63
1:A:185:HIS:HE1	1:A:187:ARG:HE	1.45	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:235:MET:CE	1:D:237:ILE:HD11	2.30	0.62
1:C:47:GLN:O	1:C:50:THR:HG22	1.99	0.62
1:A:183:SER:HB3	1:A:203:ALA:HB3	1.80	0.62
1:D:90:GLU:O	1:D:171:PHE:HA	2.00	0.62
1:B:90:GLU:O	1:B:171:PHE:HA	1.99	0.61
1:A:251:THR:HG22	1:A:359:LEU:HD21	1.81	0.61
1:A:87:ASN:O	1:A:90:GLU:OE2	2.19	0.61
1:D:35:SER:H	1:D:322:GLN:HE22	1.46	0.61
1:A:17:VAL:HG22	1:A:373:MET:HE1	1.81	0.60
1:A:259:ILE:HD13	1:A:361:VAL:HG11	1.83	0.60
1:D:154:THR:O	1:D:157:VAL:CG1	2.50	0.60
1:B:165:LEU:C	1:B:165:LEU:HD23	2.22	0.60
1:D:243:LYS:O	1:D:244:GLU:CB	2.48	0.60
1:A:100:GLN:HE22	1:A:126:PHE:H	1.49	0.60
1:B:235:MET:HE2	1:B:237:ILE:HD11	1.84	0.60
1:A:186:ARG:HB2	1:A:186:ARG:HH11	1.67	0.60
1:D:33:VAL:HG11	1:D:326:ILE:HG22	1.84	0.59
1:B:235:MET:CE	1:B:237:ILE:HD11	2.32	0.59
1:A:187:ARG:HG2	1:A:188:LEU:H	1.68	0.58
1:A:257:GLN:HG2	1:B:150:HIS:CD2	2.37	0.58
1:B:363:ARG:NH1	1:B:365:ASN:HD22	2.00	0.58
1:D:154:THR:O	1:D:157:VAL:HG12	2.03	0.58
1:B:106:VAL:HG23	1:B:109:PHE:HB2	1.86	0.57
1:B:100:GLN:HE22	1:B:126:PHE:H	1.49	0.57
1:C:331:SER:C	1:C:333:THR:H	2.08	0.57
1:D:42:VAL:HG11	1:D:167:ASN:HB2	1.86	0.57
1:B:8:VAL:HG21	1:B:74:ALA:CB	2.35	0.57
1:B:93:THR:CB	1:B:169:LEU:HD13	2.34	0.57
1:D:58:ILE:HD11	1:D:298:MET:HE3	1.87	0.57
1:C:50:THR:HG23	1:C:55:GLN:CB	2.35	0.56
1:D:33:VAL:HG11	1:D:326:ILE:CG2	2.35	0.56
1:C:42:VAL:CG2	1:C:167:ASN:HB2	2.32	0.56
1:C:296:THR:HB	1:C:300:ARG:HH21	1.70	0.56
1:A:187:ARG:HG2	1:A:188:LEU:N	2.21	0.56
1:A:296:THR:HB	1:A:300:ARG:HH21	1.71	0.55
1:C:30:ARG:HH11	1:C:30:ARG:HG3	1.72	0.55
1:D:39:VAL:O	1:D:43:LEU:HD23	2.06	0.55
1:C:53:GLU:O	1:C:57:GLN:HG3	2.07	0.55
1:D:35:SER:H	1:D:322:GLN:NE2	2.05	0.55
1:D:165:LEU:C	1:D:165:LEU:HD23	2.28	0.54
1:A:44:ALA:O	1:A:47:GLN:HB2	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:183:SER:O	1:C:185:HIS:HD2	1.91	0.54
1:C:240:PRO:HB3	1:C:245:VAL:HB	1.90	0.54
1:C:44:ALA:O	1:C:47:GLN:HB2	2.07	0.54
1:A:262:TRP:O	1:A:266:MET:HG3	2.07	0.54
1:A:243:LYS:O	1:A:244:GLU:HB2	2.07	0.54
1:A:240:PRO:HB3	1:A:245:VAL:HB	1.89	0.54
1:C:58:ILE:HD11	1:C:298:MET:CE	2.38	0.54
1:D:5:PRO:O	1:D:7:TYR:N	2.40	0.54
1:A:53:GLU:O	1:A:57:GLN:HG3	2.08	0.53
1:D:221:TYR:CE2	1:D:223:ILE:HD11	2.43	0.53
1:A:307:THR:HB	1:A:311:ASP:HB3	1.91	0.53
1:D:7:TYR:CZ	1:D:11:LEU:HD11	2.44	0.53
1:D:221:TYR:CE2	1:D:250:LEU:HD11	2.43	0.53
1:C:35:SER:H	1:C:322:GLN:NE2	1.99	0.53
1:D:235:MET:HE3	1:D:237:ILE:HD11	1.89	0.53
1:A:33:VAL:HG11	1:A:326:ILE:CG2	2.39	0.53
1:B:254:LEU:HD21	1:B:359:LEU:CD2	2.39	0.53
1:C:104:LYS:HE2	1:C:312:GLN:HB2	1.90	0.52
1:D:44:ALA:O	1:D:47:GLN:HB2	2.10	0.52
1:B:42:VAL:HG12	1:B:95:ASP:HB3	1.90	0.52
1:B:87:ASN:HD21	1:B:91:ILE:HG23	1.73	0.52
1:C:183:SER:HB3	1:C:203:ALA:HB3	1.90	0.52
1:D:87:ASN:HD21	1:D:91:ILE:CG2	2.22	0.52
1:A:92:SER:O	1:A:169:LEU:HA	2.09	0.52
1:B:291:GLU:OE2	1:B:300:ARG:NH1	2.42	0.52
1:B:313:GLU:HG2	1:B:314:PRO:HD2	1.92	0.52
1:B:330:GLU:O	1:B:331:SER:HB3	2.10	0.52
1:B:281:GLU:OE2	1:B:325:LYS:NZ	2.35	0.51
1:A:35:SER:H	1:A:322:GLN:NE2	2.00	0.51
1:D:42:VAL:HG23	1:D:43:LEU:HD22	1.92	0.51
1:B:248:SER:HA	1:B:251:THR:OG1	2.10	0.51
1:C:58:ILE:HD11	1:C:298:MET:HE2	1.91	0.51
1:A:101:ARG:HH11	1:A:101:ARG:HG3	1.76	0.51
1:C:243:LYS:O	1:C:244:GLU:CB	2.54	0.51
1:A:137:ASN:ND2	1:A:151:LEU:H	2.08	0.51
1:B:154:THR:O	1:B:157:VAL:HG13	2.10	0.51
1:C:257:GLN:HG2	1:D:150:HIS:CE1	2.46	0.51
1:D:260:SER:HB3	1:D:363:ARG:NH2	2.26	0.51
1:C:288:LYS:HB2	1:C:289:PRO:HD3	1.93	0.51
1:C:313:GLU:HG3	1:C:314:PRO:HD2	1.93	0.51
1:C:32:VAL:HG13	1:C:375:GLN:HG2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:LEU:CD1	1:B:359:LEU:HD12	2.40	0.50
1:A:185:HIS:CG	1:C:152:LEU:HD12	2.46	0.50
1:D:236:PHE:O	1:D:360:PHE:HA	2.11	0.50
1:B:148:ILE:CG2	1:B:323:LYS:HD3	2.41	0.50
1:B:168:ALA:C	1:B:169:LEU:HD22	2.30	0.50
1:C:248:SER:HA	1:C:251:THR:OG1	2.12	0.49
1:C:87:ASN:HD21	1:C:91:ILE:HG23	1.77	0.49
1:B:42:VAL:HG11	1:B:167:ASN:HB2	1.95	0.49
1:A:185:HIS:CE1	1:A:187:ARG:HE	2.27	0.49
1:B:101:ARG:HB2	1:B:123:GLN:HB3	1.95	0.49
1:B:53:GLU:O	1:B:57:GLN:HG3	2.13	0.49
1:C:99:VAL:HG22	1:C:100:GLN:N	2.28	0.49
1:C:224:LEU:HD11	1:C:273:LEU:HD13	1.94	0.49
1:C:44:ALA:HB1	1:C:66:ILE:HD13	1.94	0.49
1:A:137:ASN:HD21	1:A:151:LEU:H	1.60	0.49
1:B:191:LYS:NZ	1:B:378:GLU:OE2	2.42	0.49
1:A:296:THR:HB	1:A:300:ARG:NH2	2.28	0.48
1:C:185:HIS:HE1	1:C:187:ARG:HE	1.61	0.48
1:D:45:MET:HA	1:D:117:PHE:CZ	2.48	0.48
1:D:69:LYS:HA	1:D:69:LYS:HE3	1.94	0.48
1:B:93:THR:HB	1:B:169:LEU:HD13	1.93	0.48
1:C:300:ARG:HB3	1:C:303:GLN:HB2	1.95	0.48
1:B:95:ASP:HB2	2:B:401:EMJ:O10	2.13	0.48
1:B:210:TYR:CG	1:B:211:THR:N	2.81	0.48
1:B:8:VAL:HG21	1:B:74:ALA:HB3	1.95	0.48
1:D:247:LEU:HD21	1:D:359:LEU:HD22	1.96	0.48
1:D:263:LYS:HD3	1:D:263:LYS:O	2.13	0.48
1:B:235:MET:HA	1:B:361:VAL:O	2.13	0.48
1:C:87:ASN:ND2	1:C:91:ILE:HG23	2.28	0.48
1:B:282:THR:HG22	1:B:284:VAL:HG13	1.95	0.48
1:B:271:ARG:HD3	1:B:350:GLU:HB2	1.95	0.48
1:B:8:VAL:HG21	1:B:74:ALA:HB1	1.96	0.48
1:C:110:MET:HB2	1:C:111:PRO:HD3	1.95	0.47
1:D:16:GLY:HA3	1:D:36:PRO:HB3	1.96	0.47
1:C:55:GLN:O	1:C:59:GLN:HG3	2.13	0.47
1:C:38:GLY:O	1:C:42:VAL:HG23	2.15	0.47
1:A:258:LEU:HD13	1:B:152:LEU:HG	1.97	0.47
1:A:33:VAL:HG11	1:A:326:ILE:HG22	1.95	0.47
1:B:137:ASN:ND2	1:B:151:LEU:H	2.11	0.47
1:D:235:MET:HE2	1:D:237:ILE:HD11	1.96	0.47
1:C:364:HIS:HD2	1:C:367:THR:OG1	1.98	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:THR:HG22	1:C:109:PHE:HZ	1.80	0.47
1:D:307:THR:HA	1:D:310:SER:O	2.14	0.47
1:D:87:ASN:HD21	1:D:91:ILE:HG23	1.80	0.47
1:D:106:VAL:CG2	1:D:109:PHE:HB2	2.45	0.47
1:A:104:LYS:HE2	1:A:312:GLN:HB2	1.96	0.47
1:B:87:ASN:HD21	1:B:91:ILE:CG2	2.28	0.47
1:A:179:PHE:O	1:A:180:PRO:C	2.52	0.46
1:A:221:TYR:CZ	1:A:250:LEU:HB2	2.50	0.46
1:B:282:THR:O	1:B:323:LYS:HA	2.16	0.46
1:A:28:LYS:O	1:A:29:ASP:HB2	2.15	0.46
1:A:8:VAL:HG21	1:A:74:ALA:CB	2.45	0.46
1:B:35:SER:H	1:B:322:GLN:HE22	1.61	0.46
1:A:50:THR:HG23	1:A:55:GLN:HB2	1.97	0.46
1:A:329:ASN:HD22	1:A:331:SER:H	1.64	0.46
1:B:284:VAL:HG23	1:B:284:VAL:O	2.15	0.46
1:B:221:TYR:CD1	1:B:223:ILE:HD11	2.51	0.46
1:C:88:LYS:HB2	1:C:88:LYS:NZ	2.31	0.46
1:D:287:ARG:O	1:D:291:GLU:HG3	2.15	0.46
1:A:101:ARG:N	1:A:124:VAL:O	2.47	0.46
1:A:32:VAL:HG13	1:A:375:GLN:HG2	1.97	0.46
1:B:110:MET:HB2	1:B:111:PRO:HD3	1.98	0.46
1:B:223:ILE:HA	1:B:237:ILE:O	2.15	0.46
1:B:243:LYS:HE3	1:B:356:ARG:HB3	1.98	0.46
1:D:210:TYR:CG	1:D:211:THR:N	2.84	0.45
2:B:401:EMJ:H5	2:B:401:EMJ:O9	2.16	0.45
1:D:87:ASN:O	1:D:90:GLU:HG3	2.16	0.45
1:A:168:ALA:C	1:A:169:LEU:HD23	2.37	0.45
1:B:48:LEU:HD11	1:B:112:HIS:CE1	2.51	0.45
1:B:365:ASN:N	1:B:366:PRO:CD	2.79	0.45
1:D:247:LEU:HD11	1:D:359:LEU:CD2	2.47	0.45
1:D:21:GLN:O	1:D:25:GLN:HG3	2.16	0.45
1:B:239:ALA:HB1	1:B:356:ARG:HD2	1.98	0.45
1:B:35:SER:H	1:B:322:GLN:NE2	2.15	0.45
1:C:49:THR:HG22	1:C:109:PHE:CZ	2.52	0.45
1:A:140:VAL:HG12	1:A:148:ILE:O	2.16	0.45
1:A:52:GLY:O	1:A:56:GLN:HG2	2.17	0.45
1:D:313:GLU:HG2	1:D:314:PRO:N	2.31	0.45
1:B:287:ARG:O	1:B:291:GLU:HG3	2.17	0.45
1:C:262:TRP:O	1:C:266:MET:HG3	2.16	0.45
1:D:32:VAL:HG13	1:D:375:GLN:HG2	1.99	0.45
1:B:106:VAL:CG2	1:B:109:PHE:HB2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:LEU:HD23	1:C:165:LEU:C	2.37	0.44
1:C:28:LYS:HG2	1:C:29:ASP:OD1	2.17	0.44
1:A:281:GLU:HB2	1:A:325:LYS:HG2	1.99	0.44
1:C:271:ARG:HE	1:C:350:GLU:HB2	1.83	0.44
1:D:206:ASN:OD1	1:D:207:LYS:N	2.43	0.44
1:A:185:HIS:CE1	1:A:187:ARG:HH21	2.35	0.44
1:A:18:ARG:NH2	1:B:130:GLU:OE1	2.51	0.44
1:A:185:HIS:CD2	1:C:152:LEU:HD12	2.53	0.44
1:D:167:ASN:O	1:D:322:GLN:HA	2.17	0.44
1:A:243:LYS:O	1:A:244:GLU:CB	2.65	0.44
1:A:45:MET:HA	1:A:117:PHE:CZ	2.52	0.44
1:C:167:ASN:O	1:C:322:GLN:HA	2.17	0.44
1:D:196:THR:O	1:D:197:VAL:HG13	2.18	0.44
1:A:305:ASP:OD1	1:A:305:ASP:C	2.55	0.44
1:B:202:MET:HG3	1:B:277:LYS:HG2	2.00	0.44
1:D:221:TYR:CE2	1:D:223:ILE:CD1	3.01	0.44
1:A:185:HIS:HE1	1:A:187:ARG:NE	2.14	0.44
1:C:140:VAL:HG12	1:C:148:ILE:O	2.18	0.44
1:D:233:LEU:HD23	1:D:364:HIS:HA	2.00	0.44
1:A:257:GLN:HG2	1:B:150:HIS:NE2	2.32	0.44
1:D:106:VAL:HG23	1:D:109:PHE:HB2	2.00	0.44
1:A:300:ARG:HG2	1:A:303:GLN:NE2	2.33	0.43
1:A:364:HIS:HD2	1:A:367:THR:OG1	2.01	0.43
1:B:87:ASN:O	1:B:90:GLU:OE2	2.36	0.43
1:C:342:ILE:HG22	1:C:343:VAL:N	2.33	0.43
1:A:363:ARG:HH11	1:A:365:ASN:HD22	1.64	0.43
1:B:7:TYR:CZ	1:B:11:LEU:HD11	2.54	0.43
1:B:251:THR:HG22	1:B:359:LEU:HD21	2.00	0.43
1:D:38:GLY:HA3	1:D:167:ASN:HD21	1.83	0.43
1:D:8:VAL:HG21	1:D:74:ALA:HB3	1.99	0.43
1:A:110:MET:HB2	1:A:111:PRO:HD3	2.00	0.43
2:B:401:EMJ:C13	2:B:401:EMJ:O9	2.66	0.43
1:A:254:LEU:HD13	1:A:258:LEU:HD23	2.01	0.43
1:B:236:PHE:O	1:B:360:PHE:HA	2.19	0.43
1:B:5:PRO:N	1:B:8:VAL:HG23	2.34	0.43
1:D:58:ILE:HD11	1:D:298:MET:CE	2.49	0.43
1:B:222:ASP:O	1:B:238:ALA:HA	2.19	0.42
1:B:8:VAL:HG11	1:B:75:LEU:HG	2.01	0.42
1:A:258:LEU:O	1:A:261:HIS:HB3	2.19	0.42
1:C:190:HIS:O	1:C:357:PRO:HD3	2.19	0.42
1:A:72:ALA:N	1:A:73:PRO:CD	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:ASN:HD21	1:B:332:GLY:N	2.17	0.42
1:B:363:ARG:HH11	1:B:365:ASN:HD22	1.63	0.42
1:D:46:LEU:HD23	1:D:298:MET:HE1	2.00	0.42
1:C:157:VAL:HG23	1:C:157:VAL:O	2.19	0.42
1:A:165:LEU:HD23	1:A:165:LEU:C	2.40	0.42
1:A:43:LEU:HA	1:A:43:LEU:HD13	1.84	0.42
1:A:93:THR:O	1:A:94:THR:HB	2.20	0.42
1:D:278:PHE:CZ	1:D:328:VAL:HG21	2.55	0.42
1:D:224:LEU:HD11	1:D:352:ILE:HG21	2.01	0.42
1:A:311:ASP:N	1:A:311:ASP:OD1	2.51	0.42
1:A:87:ASN:HD21	1:A:91:ILE:HG23	1.84	0.42
1:C:72:ALA:N	1:C:73:PRO:CD	2.83	0.42
1:D:48:LEU:HD21	1:D:116:LEU:HD22	2.02	0.42
1:B:307:THR:HA	1:B:310:SER:O	2.20	0.42
1:C:187:ARG:CG	1:C:188:LEU:N	2.83	0.42
1:A:50:THR:HG23	1:A:55:GLN:CB	2.50	0.41
1:B:148:ILE:HG23	1:B:323:LYS:HD3	2.02	0.41
1:D:109:PHE:CD1	1:D:109:PHE:C	2.93	0.41
1:A:288:LYS:HB2	1:A:289:PRO:HD3	2.03	0.41
1:D:282:THR:O	1:D:323:LYS:HA	2.19	0.41
1:B:207:LYS:HA	1:B:269:LEU:O	2.20	0.41
1:A:140:VAL:HG11	1:A:148:ILE:HG22	2.02	0.41
1:C:185:HIS:HE1	1:C:187:ARG:NE	2.19	0.41
1:C:331:SER:C	1:C:333:THR:N	2.73	0.41
1:D:140:VAL:HG12	1:D:148:ILE:O	2.20	0.41
1:A:215:THR:HG22	1:A:253:ILE:CD1	2.51	0.41
1:B:16:GLY:HA3	1:B:36:PRO:HB3	2.02	0.41
1:B:191:LYS:NZ	1:B:378:GLU:CD	2.73	0.41
1:C:292:ASN:O	1:D:159:GLN:NE2	2.54	0.41
1:B:247:LEU:HD11	1:B:359:LEU:CD1	2.45	0.41
1:C:86:TRP:CE3	1:C:87:ASN:HA	2.56	0.41
1:D:204:GLN:HG3	1:D:208:PHE:HZ	1.86	0.41
1:D:238:ALA:HB1	1:D:250:LEU:HD21	2.03	0.41
1:D:33:VAL:HG12	1:D:360:PHE:HZ	1.85	0.41
1:B:122:LYS:NZ	2:B:401:EMJ:H11	2.35	0.41
1:D:229:HIS:O	1:D:229:HIS:CG	2.73	0.41
1:B:44:ALA:O	1:B:47:GLN:HB2	2.21	0.41
1:D:235:MET:HA	1:D:361:VAL:O	2.21	0.41
1:B:167:ASN:O	1:B:322:GLN:HA	2.21	0.41
1:B:139:TRP:HZ2	2:B:401:EMJ:H5	1.86	0.41
1:B:119:SER:HB3	2:B:401:EMJ:H6	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:154:THR:O	1:D:157:VAL:HG13	2.20	0.41
1:D:126:PHE:O	1:D:159:GLN:HA	2.21	0.41
1:B:119:SER:OG	2:B:401:EMJ:H4	2.21	0.40
1:B:201:MET:HA	1:B:276:PRO:HA	2.03	0.40
1:A:90:GLU:O	1:A:171:PHE:HA	2.21	0.40
1:C:100:GLN:HE22	1:C:126:PHE:H	1.69	0.40
1:C:310:SER:HB3	1:C:315:LEU:HD12	2.03	0.40
1:A:35:SER:N	1:A:322:GLN:HE22	2.01	0.40
1:C:329:ASN:HD22	1:C:329:ASN:C	2.24	0.40
1:A:162:ARG:HD2	1:A:313:GLU:OE1	2.22	0.40
1:B:33:VAL:HG21	1:B:280:LEU:HB2	2.03	0.40
1:C:224:LEU:CD1	1:C:273:LEU:HD13	2.52	0.40
1:D:87:ASN:ND2	1:D:91:ILE:HG22	2.36	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	363/375 (97%)	340 (94%)	16 (4%)	7 (2%)	8	9
1	B	356/375 (95%)	340 (96%)	12 (3%)	4 (1%)	14	19
1	C	364/375 (97%)	344 (94%)	12 (3%)	8 (2%)	6	7
1	D	356/375 (95%)	338 (95%)	14 (4%)	4 (1%)	14	19
All	All	1439/1500 (96%)	1362 (95%)	54 (4%)	23 (2%)	9	12

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	193	ASP
1	B	87	ASN

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Mol	Chain	Res	Type
1	B	331	SER
1	A	87	ASN
1	A	129	VAL
1	B	88	LYS
1	C	101	ARG
1	D	6	SER
1	D	244	GLU
1	A	88	LYS
1	A	331	SER
1	B	244	GLU
1	C	87	ASN
1	C	88	LYS
1	C	333	THR
1	C	102	ASP
1	C	244	GLU
1	A	244	GLU
1	D	87	ASN
1	D	314	PRO
1	A	155	GLY
1	C	349	PRO
1	C	129	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	322/329 (98%)	300 (93%)	22 (7%)	16	24
1	B	317/329 (96%)	298 (94%)	19 (6%)	19	30
1	C	323/329 (98%)	310 (96%)	13 (4%)	31	48
1	D	317/329 (96%)	293 (92%)	24 (8%)	13	20
All	All	1279/1316 (97%)	1201 (94%)	78 (6%)	18	29

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	32	VAL
1	A	50	THR
1	A	53	GLU
1	A	102	ASP
1	A	130	GLU
1	A	149	SER
1	A	158	ASP
1	A	165	LEU
1	A	169	LEU
1	A	182	SER
1	A	186	ARG
1	A	215	THR
1	A	242	GLU
1	A	248	SER
1	A	254	LEU
1	A	255	SER
1	A	271	ARG
1	A	284	VAL
1	A	312	GLN
1	A	329	ASN
1	A	346	ARG
1	B	8	VAL
1	B	21	GLN
1	B	32	VAL
1	B	69	LYS
1	B	91	ILE
1	B	101	ARG
1	B	107	GLN
1	B	145	LYS
1	B	152	LEU
1	B	157	VAL
1	B	159	GLN
1	B	165	LEU
1	B	169	LEU
1	B	196	THR
1	B	254	LEU
1	B	257	GLN
1	B	321	LEU
1	B	329	ASN
1	B	356	ARG
1	C	23	VAL
1	C	29	ASP

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Mol	Chain	Res	Type
1	C	32	VAL
1	C	43	LEU
1	C	50	THR
1	C	80	LYS
1	C	88	LYS
1	C	169	LEU
1	C	212	GLU
1	C	265	ASN
1	C	284	VAL
1	C	329	ASN
1	C	346	ARG
1	D	8	VAL
1	D	32	VAL
1	D	67	ASP
1	D	69	LYS
1	D	101	ARG
1	D	127	SER
1	D	152	LEU
1	D	157	VAL
1	D	159	GLN
1	D	165	LEU
1	D	169	LEU
1	D	172	ASN
1	D	181	ASP
1	D	186	ARG
1	D	223	ILE
1	D	224	LEU
1	D	257	GLN
1	D	263	LYS
1	D	284	VAL
1	D	298	MET
1	D	300	ARG
1	D	308	SER
1	D	321	LEU
1	D	325	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (55) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	59	GLN
1	A	100	GLN

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Mol	Chain	Res	Type
1	A	137	ASN
1	A	167	ASN
1	A	172	ASN
1	A	185	HIS
1	A	190	HIS
1	A	292	ASN
1	A	301	GLN
1	A	303	GLN
1	A	312	GLN
1	A	322	GLN
1	A	329	ASN
1	A	364	HIS
1	A	365	ASN
1	B	21	GLN
1	B	59	GLN
1	B	100	GLN
1	B	107	GLN
1	B	112	HIS
1	B	137	ASN
1	B	143	HIS
1	B	167	ASN
1	B	292	ASN
1	B	322	GLN
1	B	329	ASN
1	B	364	HIS
1	B	365	ASN
1	C	21	GLN
1	C	25	GLN
1	C	100	GLN
1	C	112	HIS
1	C	137	ASN
1	C	159	GLN
1	C	167	ASN
1	C	172	ASN
1	C	185	HIS
1	C	204	GLN
1	C	265	ASN
1	C	292	ASN
1	C	322	GLN
1	C	329	ASN
1	C	364	HIS
1	C	365	ASN

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Mol	Chain	Res	Type
1	D	25	GLN
1	D	167	ASN
1	D	172	ASN
1	D	257	GLN
1	D	261	HIS
1	D	292	ASN
1	D	303	GLN
1	D	322	GLN
1	D	364	HIS
1	D	365	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	EMJ	B	401	-	21,21,21	1.38	3 (14%)	24,26,26	2.16	3 (12%)

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	EMJ	B	401	-	-	5/11/31/31	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	EMJ	C6-C7	3.31	1.40	1.35
2	B	401	EMJ	C4-C5	-2.89	1.39	1.45
2	B	401	EMJ	C6-C5	-2.02	1.39	1.44

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	EMJ	C12-C3-C4	-6.21	116.05	123.08
2	B	401	EMJ	C12-C3-C2	5.67	126.28	117.00
2	B	401	EMJ	O9-C2-C7	-3.96	114.33	120.14

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	EMJ	C13-C12-C3-C2
2	B	401	EMJ	C13-C12-C3-C4
2	B	401	EMJ	C14-C15-C16-C17
2	B	401	EMJ	C12-C13-C14-C15
2	B	401	EMJ	C16-C17-C18-C19

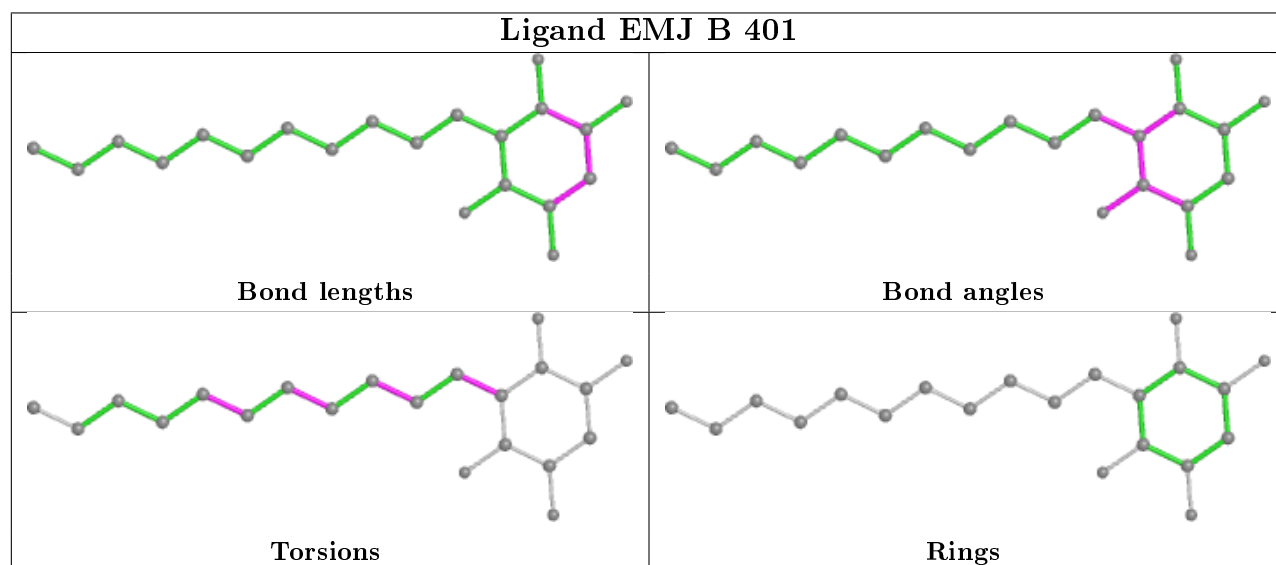
There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	EMJ	7	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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 ⓘ

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	367/375 (97%)	-0.30	0 100 100	15, 31, 61, 97	0
1	B	360/375 (96%)	-0.18	5 (1%) 75 73	18, 34, 65, 102	0
1	C	368/375 (98%)	-0.26	3 (0%) 86 84	16, 31, 62, 95	0
1	D	360/375 (96%)	-0.15	5 (1%) 75 73	17, 35, 67, 98	0
All	All	1455/1500 (97%)	-0.22	13 (0%) 84 82	15, 33, 64, 102	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	7	TYR	3.9
1	C	334	VAL	3.9
1	B	312	GLN	3.1
1	C	104	LYS	3.0
1	D	312	GLN	2.6
1	B	88	LYS	2.5
1	D	102	ASP	2.5
1	B	313	GLU	2.4
1	D	231	ASP	2.3
1	D	87	ASN	2.3
1	B	87	ASN	2.2
1	B	86	TRP	2.2
1	D	313	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

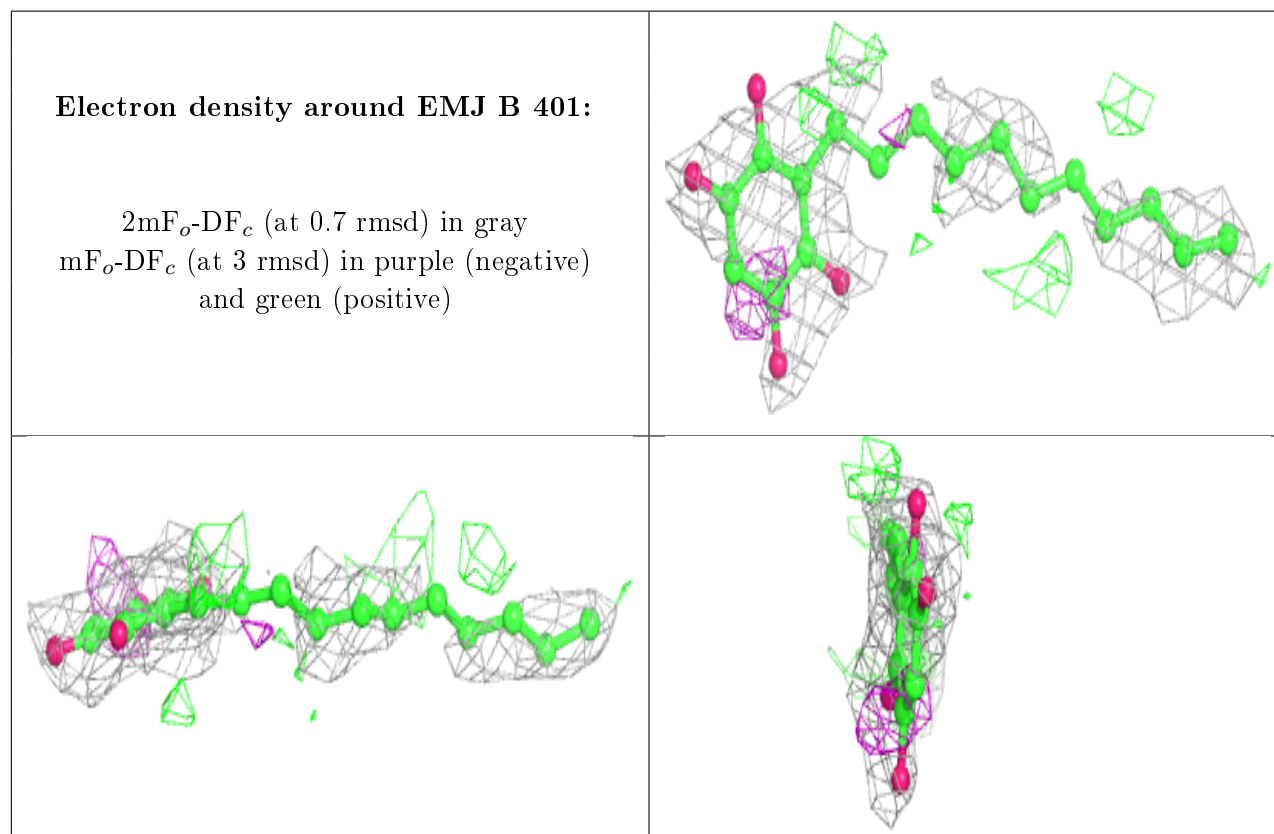
There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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(\AA^2)	Q<0.9
2	EMJ	B	401	21/21	0.69	0.33	57,63,74,75	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.