



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:16 AM UTC

PDB ID : 1ME6 / pdb_00001me6
Title : CRYSTAL STRUCTURE OF PLASMEPSIN II, AN ASPARTYL PROTEASE FROM PLASMODIUM FALCIPARUM, IN COMPLEX WITH A STATINE-BASED INHIBITOR
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Deposited on : 2002-08-08
Resolution : 2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtrriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

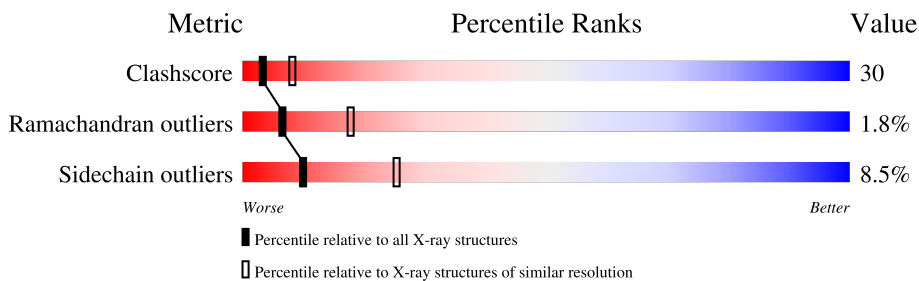
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	329	
1	B	329	

2 Entry composition [i](#)

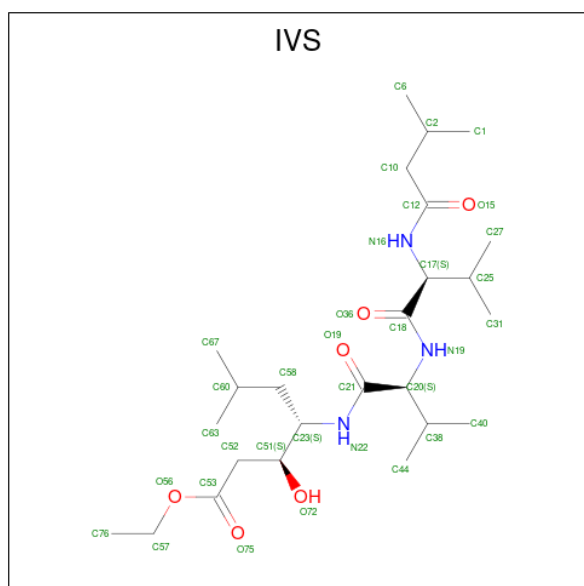
There are 3 unique types of molecules in this entry. The entry contains 5402 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Plasmepsin II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2607	1689	404	503	11			
1	B	329	Total	C	N	O	S	0	0	0
			2607	1689	404	503	11			

- Molecule 2 is 3-HYDROXY-6-METHYL-4-(3-METHYL-2-(3-METHYL-2-(3-METHYL-BUTYRYLAMINO)-BUTYRYLAMINO)-BUTYRYLAMINO)-HEPTANOIC ACID ETHYL ESTER (CCD ID: IVS) (formula: $C_{25}H_{47}N_3O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			34	25	3	6		
2	B	1	Total	C	N	O	0	0
			34	25	3	6		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	82	Total 82	O 82	0	0
3	B	38	Total 38	O 38	0	0

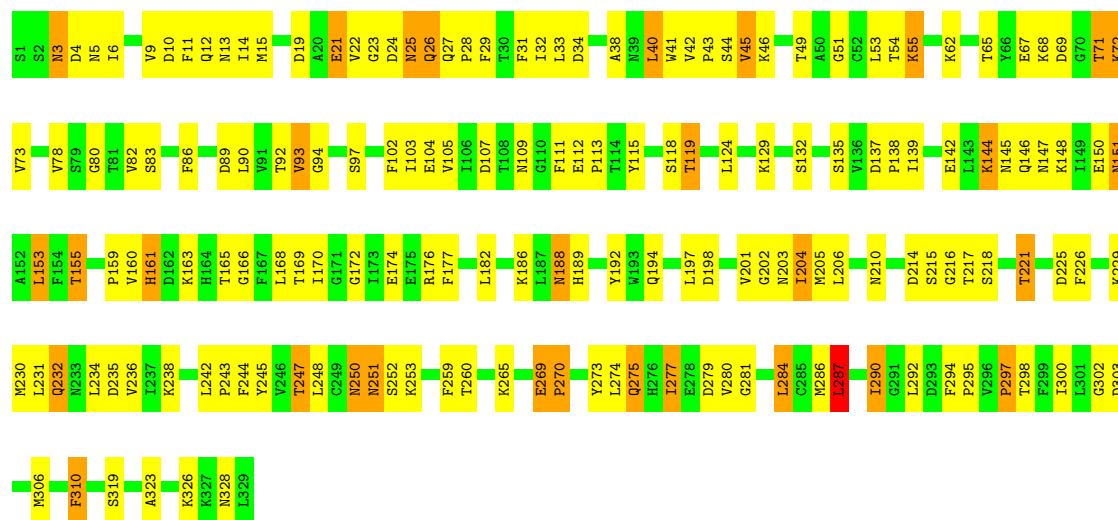
3 Residue-property plots

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: Plasmepsin II

Chain A: 



T311	V312	Y315	H318	A323	K326	K327	N328	L329
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4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	141.93Å 141.93Å 97.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.70	Depositor
% Data completeness (in resolution range)	89.8 (50.00-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.214 , 0.271	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5402	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.50	0/2674	1.05	11/3638 (0.3%)
1	B	0.47	0/2674	1.06	19/3638 (0.5%)
All	All	0.49	0/5348	1.06	30/7276 (0.4%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	248	LEU	N-CA-C	-8.72	95.89	109.76
1	B	280	VAL	N-CA-C	-8.23	105.12	111.62
1	B	27	GLN	CA-C-N	8.04	128.19	120.31
1	B	27	GLN	C-N-CA	8.04	128.19	120.31
1	A	287	LEU	N-CA-C	7.60	121.81	112.23
1	A	32	ILE	N-CA-C	-7.51	96.25	107.37
1	B	14	ILE	N-CA-C	-6.94	105.59	111.56
1	A	177	PHE	N-CA-C	6.61	118.56	111.36
1	B	263	ASN	N-CA-C	6.52	120.45	112.23
1	A	55	LYS	N-CA-C	6.43	119.91	110.59
1	A	25	ASN	N-CA-C	-6.42	101.02	110.52
1	B	248	LEU	N-CA-C	-6.42	99.56	109.76
1	A	192	TYR	N-CA-C	-6.34	101.08	110.46
1	B	32	ILE	N-CA-C	-6.32	98.64	107.80
1	B	224	THR	N-CA-C	6.29	118.66	111.11
1	B	35	THR	N-CA-C	-6.28	105.58	113.18
1	B	146	GLN	N-CA-C	-6.19	105.83	113.19
1	B	281	GLY	CA-C-N	6.01	127.36	119.84
1	B	281	GLY	C-N-CA	6.01	127.36	119.84
1	B	2	SER	N-CA-C	-5.76	106.25	113.28
1	A	203	ASN	N-CA-C	-5.73	106.29	113.28
1	B	193	TRP	N-CA-C	-5.62	100.46	109.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	45	VAL	N-CA-C	-5.59	105.46	113.07
1	B	167	PHE	N-CA-C	5.38	118.20	109.06
1	B	192	TYR	N-CA-C	-5.29	102.91	110.59
1	A	204	ILE	N-CA-C	5.21	116.08	108.58
1	B	90	LEU	N-CA-C	5.17	116.89	108.63
1	B	310	PHE	N-CA-C	-5.11	102.08	109.59
1	A	243	PRO	N-CA-C	-5.08	104.77	113.70
1	A	310	PHE	N-CA-C	-5.04	102.92	110.23

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	2607	0	2536	158	0
1	B	2607	0	2536	160	0
2	A	34	0	47	2	0
2	B	34	0	47	1	0
3	A	82	0	0	10	0
3	B	38	0	0	4	0
All	All	5402	0	5166	318	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 30.

All (318) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:HIS:HA	1:B:285:CYS:HB3	1.43	1.00
1:A:260:THR:HG22	1:A:265:LYS:HB2	1.44	0.97
1:A:151:ASN:HD22	1:A:151:ASN:H	1.02	0.96
1:A:69:ASP:OD1	1:A:71:THR:HG23	1.66	0.96
1:A:45:VAL:HG12	1:A:46:LYS:HD2	1.46	0.95
1:A:90:LEU:HD11	1:A:97:SER:HB3	1.50	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:GLN:HA	3:A:873:HOH:O	1.70	0.91
1:B:201:VAL:HG11	1:B:230:MET:HE3	1.53	0.90
1:B:44:SER:HB2	1:B:104:GLU:HG2	1.55	0.88
1:A:151:ASN:H	1:A:151:ASN:ND2	1.68	0.84
1:B:163:LYS:HG2	1:B:327:LYS:HE3	1.57	0.84
1:B:221:THR:HB	1:B:300:ILE:HB	1.61	0.82
1:A:151:ASN:HD22	1:A:151:ASN:N	1.76	0.81
1:A:182:LEU:HD23	1:A:323:ALA:HB2	1.61	0.81
1:B:281:GLY:HA3	1:B:284:LEU:HD22	1.64	0.79
1:A:251:ASN:ND2	1:A:253:LYS:H	1.81	0.79
1:B:142:GLU:O	1:B:146:GLN:HG2	1.83	0.78
1:B:144:LYS:HA	1:B:149:ILE:HG22	1.66	0.77
1:A:198:ASP:HB3	1:A:205:MET:HE1	1.66	0.77
1:B:206:LEU:HD23	1:B:299:PHE:HE1	1.49	0.76
1:B:277:ILE:CG2	1:B:280:VAL:HB	2.15	0.76
1:A:221:THR:HB	1:A:300:ILE:HB	1.68	0.76
1:A:151:ASN:ND2	1:A:151:ASN:N	2.29	0.74
1:B:251:ASN:HB3	1:B:254:LEU:HG	1.70	0.73
1:B:222:VAL:HG23	1:B:299:PHE:CE2	2.23	0.73
1:A:21:GLU:HB2	1:A:26:GLN:O	1.90	0.72
1:B:41:TRP:CE3	1:B:105:VAL:HG21	2.24	0.72
1:B:25:ASN:O	1:B:26:GLN:HB3	1.87	0.72
1:B:163:LYS:NZ	1:B:327:LYS:HG3	2.05	0.71
1:A:235:ASP:N	3:A:924:HOH:O	2.15	0.71
1:B:152:ALA:O	1:B:315:TYR:HB2	1.91	0.71
1:A:69:ASP:HB3	1:A:86:PHE:O	1.92	0.69
1:A:236:VAL:HG22	1:A:247:THR:CG2	2.22	0.69
1:A:6:ILE:CD1	1:A:93:VAL:HG12	2.23	0.69
1:A:221:THR:HG23	1:A:292:LEU:HB3	1.75	0.69
1:B:149:ILE:HG12	1:B:170:ILE:HG22	1.74	0.69
1:A:260:THR:HG22	1:A:265:LYS:CB	2.23	0.68
1:B:23:GLY:HA2	1:B:89:ASP:OD1	1.92	0.68
1:A:174:GLU:OE1	1:A:176:ARG:HD3	1.94	0.67
1:B:21:GLU:HG2	1:B:92:THR:HB	1.76	0.67
1:A:129:LYS:HD2	1:A:137:ASP:OD1	1.94	0.67
1:B:49:THR:HG21	1:B:108:THR:OG1	1.95	0.67
1:B:276:HIS:HA	1:B:285:CYS:CB	2.24	0.67
1:B:129:LYS:HE2	1:B:135:SER:HB2	1.76	0.67
1:A:3:ASN:HD22	1:A:172:GLY:HA3	1.60	0.66
1:B:146:GLN:HB2	1:B:148:LYS:HG3	1.76	0.66
1:B:223:PRO:HG2	1:B:226:PHE:HB2	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:ASP:OD2	1:A:13:ASN:HA	1.95	0.66
1:A:236:VAL:HG22	1:A:247:THR:HG23	1.77	0.66
1:A:188:ASN:C	1:A:188:ASN:HD22	2.04	0.66
1:B:247:THR:HG22	1:B:248:LEU:O	1.96	0.66
1:A:279:ASP:HB2	3:A:823:HOH:O	1.97	0.65
1:A:90:LEU:CD1	1:A:97:SER:HB3	2.25	0.65
1:A:21:GLU:HG2	1:A:92:THR:HB	1.80	0.64
1:A:160:VAL:HG11	1:A:163:LYS:HE2	1.79	0.64
1:A:153:LEU:C	1:A:153:LEU:HD12	2.22	0.64
1:B:163:LYS:HZ1	1:B:327:LYS:HG3	1.62	0.64
1:B:258:GLU:HG2	1:B:267:THR:HG22	1.77	0.64
1:A:251:ASN:C	1:A:251:ASN:HD22	2.04	0.64
1:A:182:LEU:HD23	1:A:323:ALA:CB	2.27	0.64
1:A:277:ILE:HG23	1:A:277:ILE:O	1.97	0.63
1:B:153:LEU:C	1:B:153:LEU:HD12	2.24	0.63
1:B:258:GLU:OE2	1:B:265:LYS:NZ	2.29	0.63
1:A:159:PRO:HB3	1:A:166:GLY:N	2.14	0.62
1:A:277:ILE:HG22	1:A:284:LEU:HB3	1.81	0.62
1:B:277:ILE:HG22	1:B:280:VAL:HB	1.81	0.61
1:A:142:GLU:O	1:A:146:GLN:HG3	2.01	0.61
1:B:277:ILE:HG22	1:B:277:ILE:O	1.99	0.61
1:B:25:ASN:HB3	1:B:27:GLN:HG3	1.81	0.61
1:B:86:PHE:HD1	1:B:102:PHE:O	1.84	0.61
1:B:247:THR:HG22	1:B:248:LEU:N	2.15	0.60
1:A:146:GLN:HB2	1:A:148:LYS:HG3	1.84	0.60
1:A:201:VAL:HG23	1:A:226:PHE:CZ	2.36	0.60
1:B:150:GLU:HG2	1:B:171:GLY:O	2.01	0.60
1:A:161:HIS:HB3	3:A:842:HOH:O	2.02	0.60
1:B:144:LYS:CA	1:B:149:ILE:HG22	2.31	0.60
1:A:3:ASN:ND2	1:A:172:GLY:HA3	2.16	0.59
1:B:84:GLY:HA2	1:B:106:ILE:HG12	1.83	0.59
1:B:222:VAL:HG21	1:B:226:PHE:CD1	2.38	0.59
1:B:281:GLY:HA3	1:B:284:LEU:CD2	2.33	0.59
1:B:277:ILE:HG23	1:B:280:VAL:HB	1.85	0.59
1:A:83:SER:HB2	1:A:107:ASP:HB3	1.85	0.58
1:A:6:ILE:HD13	1:A:93:VAL:HG12	1.85	0.58
1:A:197:LEU:HD13	1:A:259:PHE:HB3	1.86	0.58
1:B:234:LEU:HG	1:B:236:VAL:HG13	1.85	0.58
1:A:6:ILE:HD11	1:A:93:VAL:HG12	1.86	0.58
1:B:39:ASN:HA	3:B:830:HOH:O	2.03	0.58
1:B:149:ILE:HG13	1:B:171:GLY:HA2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:VAL:HG23	1:A:226:PHE:HZ	1.69	0.57
1:A:277:ILE:HG13	1:A:280:VAL:CG2	2.35	0.57
1:A:82:VAL:HG13	1:A:105:VAL:CG1	2.34	0.57
1:A:188:ASN:ND2	1:A:194:GLN:HG3	2.20	0.57
1:B:15:MET:HE3	1:B:16:PHE:O	2.03	0.57
1:B:34:ASP:OD1	1:B:216:GLY:HA3	2.04	0.57
1:B:149:ILE:HG13	1:B:171:GLY:CA	2.35	0.57
1:A:273:TYR:O	1:A:287:LEU:O	2.22	0.57
1:B:246:VAL:HG22	1:B:286:MET:CE	2.35	0.57
1:B:155:THR:O	1:B:168:LEU:HD12	2.05	0.57
1:A:155:THR:HG22	1:A:169:THR:HB	1.86	0.57
1:B:203:ASN:O	1:B:203:ASN:ND2	2.38	0.56
1:A:251:ASN:HD22	1:A:252:SER:N	2.03	0.56
1:A:22:VAL:O	1:A:27:GLN:HB2	2.05	0.56
1:A:188:ASN:HD21	1:A:194:GLN:HE21	1.53	0.56
1:A:281:GLY:O	1:A:284:LEU:HB2	2.05	0.56
1:B:21:GLU:CD	1:B:21:GLU:N	2.64	0.56
1:A:90:LEU:HD11	1:A:97:SER:CB	2.29	0.55
1:B:69:ASP:O	1:B:71:THR:N	2.39	0.55
1:B:201:VAL:HG11	1:B:230:MET:CE	2.33	0.55
1:B:146:GLN:HB3	1:B:148:LYS:NZ	2.21	0.55
1:B:163:LYS:CG	1:B:327:LYS:HE3	2.33	0.55
1:B:260:THR:HG22	1:B:265:LYS:HG3	1.88	0.55
1:A:251:ASN:HD22	1:A:253:LYS:H	1.54	0.55
1:A:82:VAL:HG11	1:A:105:VAL:HG21	1.88	0.55
1:B:81:THR:HG22	1:B:82:VAL:N	2.21	0.55
1:A:24:ASP:HB3	1:A:65:THR:OG1	2.07	0.55
1:A:41:TRP:CE3	1:A:105:VAL:HG21	2.42	0.54
1:A:201:VAL:HG12	1:A:202:GLY:N	2.23	0.54
1:B:206:LEU:HD23	1:B:299:PHE:CE1	2.38	0.54
1:A:53:LEU:C	1:A:53:LEU:HD23	2.33	0.54
1:A:55:LYS:NZ	1:A:118:SER:O	2.40	0.54
1:B:198:ASP:O	1:B:259:PHE:HA	2.07	0.54
1:A:155:THR:CG2	1:A:169:THR:HB	2.38	0.53
1:A:214:ASP:O	1:A:302:GLY:HA2	2.07	0.53
1:A:277:ILE:HG23	1:A:280:VAL:HB	1.89	0.53
1:A:72:LYS:H	1:A:72:LYS:HD2	1.73	0.53
1:B:277:ILE:HG13	1:B:286:MET:HE3	1.91	0.53
1:B:145:ASN:C	1:B:147:ASN:H	2.16	0.53
1:A:144:LYS:HE3	1:A:150:GLU:O	2.09	0.53
1:A:188:ASN:HD21	1:A:194:GLN:HG3	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:ASP:HB3	1:B:86:PHE:O	2.09	0.53
1:A:5:ASN:C	1:A:6:ILE:HG13	2.34	0.53
1:A:245:TYR:O	1:A:286:MET:HA	2.09	0.52
1:B:247:THR:HG21	1:B:254:LEU:HD21	1.92	0.52
1:B:16:PHE:CE2	1:B:158:LEU:HD22	2.45	0.52
1:B:66:TYR:CG	1:B:67:GLU:N	2.78	0.52
1:A:25:ASN:ND2	1:A:27:GLN:HG3	2.25	0.52
1:A:277:ILE:HG13	1:A:280:VAL:HG21	1.91	0.52
1:A:23:GLY:HA2	1:A:89:ASP:CG	2.35	0.51
1:A:145:ASN:C	1:A:147:ASN:H	2.18	0.51
1:A:186:LYS:HD3	1:A:319:SER:HB3	1.91	0.51
1:B:43:PRO:HG2	1:B:56:HIS:O	2.11	0.51
1:B:230:MET:HE2	1:B:255:PRO:HG3	1.92	0.51
1:B:269:GLU:HB3	1:B:270:PRO:HD2	1.93	0.51
1:B:277:ILE:CG2	1:B:277:ILE:O	2.59	0.50
1:B:26:GLN:OE1	1:B:26:GLN:O	2.29	0.50
1:B:251:ASN:HB3	1:B:254:LEU:CG	2.40	0.50
1:A:23:GLY:HA2	1:A:89:ASP:OD1	2.12	0.50
1:A:201:VAL:HG21	1:A:226:PHE:HE1	1.75	0.50
1:A:231:LEU:HD23	1:A:234:LEU:HD12	1.93	0.50
1:A:269:GLU:HB2	1:A:270:PRO:HD2	1.94	0.50
1:B:188:ASN:OD1	1:B:196:THR:OG1	2.26	0.50
1:B:222:VAL:HG22	1:B:226:PHE:HB3	1.92	0.50
1:A:45:VAL:CG1	1:A:46:LYS:HD2	2.32	0.50
1:A:160:VAL:HB	1:A:163:LYS:HB2	1.93	0.50
1:B:25:ASN:HB2	1:B:62:LYS:O	2.12	0.50
1:A:9:VAL:HG11	3:A:891:HOH:O	2.11	0.49
1:A:65:THR:HB	1:A:89:ASP:OD1	2.12	0.49
1:B:97:SER:O	1:B:98:LEU:HB3	2.11	0.49
1:A:45:VAL:HG12	1:A:46:LYS:CD	2.33	0.49
1:A:238:LYS:HE3	1:A:242:LEU:O	2.13	0.49
1:B:70:GLY:HA2	1:B:85:PHE:CD2	2.47	0.49
1:B:273:TYR:HA	1:B:304:PRO:HB3	1.94	0.49
1:B:248:LEU:HD12	3:B:911:HOH:O	2.12	0.49
1:B:42:VAL:CG2	1:B:43:PRO:HD2	2.42	0.49
1:B:55:LYS:HG3	1:B:115:TYR:CZ	2.48	0.49
1:A:251:ASN:ND2	1:A:251:ASN:C	2.71	0.48
1:B:13:ASN:CG	1:B:161:HIS:HB2	2.38	0.48
1:A:188:ASN:HD22	1:A:189:HIS:N	2.11	0.48
1:A:275:GLN:HG3	3:A:918:HOH:O	2.12	0.48
1:B:24:ASP:HB2	1:B:65:THR:OG1	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:THR:HG22	1:B:51:GLY:H	1.79	0.48
1:B:215:SER:HB3	1:B:306:MET:HE1	1.95	0.48
1:B:327:LYS:HD3	1:B:327:LYS:O	2.14	0.48
1:A:229:LYS:O	1:A:232:GLN:HG2	2.14	0.47
1:B:153:LEU:HD12	1:B:153:LEU:O	2.14	0.47
1:B:62:LYS:NZ	1:B:62:LYS:HB3	2.29	0.47
1:B:246:VAL:HG22	1:B:286:MET:HE1	1.95	0.47
1:A:277:ILE:HD12	1:A:277:ILE:HA	1.53	0.47
1:B:41:TRP:HE1	1:B:123:ILE:HD12	1.79	0.47
1:A:124:LEU:HD21	1:A:139:ILE:HD12	1.97	0.47
1:B:40:LEU:O	1:B:102:PHE:HB2	2.15	0.47
1:A:43:PRO:HG3	1:A:115:TYR:OH	2.15	0.47
1:A:83:SER:O	1:A:105:VAL:HG13	2.15	0.47
1:A:287:LEU:HD23	1:A:287:LEU:N	2.29	0.47
1:B:69:ASP:C	1:B:71:THR:H	2.22	0.47
1:A:15:MET:HE2	3:A:804:HOH:O	2.15	0.47
1:B:151:ASN:HB2	1:B:171:GLY:O	2.15	0.47
1:B:129:LYS:HD2	1:B:137:ASP:OD1	2.15	0.46
1:B:271:GLU:CD	1:B:271:GLU:H	2.23	0.46
1:A:4:ASP:HB3	1:A:170:ILE:HB	1.96	0.46
1:A:80:GLY:HA3	1:A:111:PHE:HD1	1.79	0.46
1:A:277:ILE:O	1:A:277:ILE:CG2	2.62	0.46
1:B:179:GLU:HB3	1:B:326:LYS:HZ3	1.79	0.46
1:A:150:GLU:HB2	1:A:151:ASN:ND2	2.29	0.46
1:A:225:ASP:O	1:A:229:LYS:HG3	2.15	0.46
1:A:146:GLN:OE1	1:A:148:LYS:NZ	2.36	0.46
1:B:60:SER:HB2	1:B:66:TYR:CE1	2.49	0.46
1:B:70:GLY:HA2	1:B:85:PHE:HD2	1.81	0.46
1:B:128:TRP:CE3	1:B:190:ASP:HB3	2.51	0.46
1:B:145:ASN:C	1:B:147:ASN:N	2.72	0.46
1:A:82:VAL:CG1	1:A:105:VAL:CG2	2.94	0.46
1:A:236:VAL:HG22	1:A:247:THR:HG21	1.95	0.46
1:B:48:THR:O	1:B:49:THR:C	2.59	0.46
1:A:34:ASP:OD1	1:A:216:GLY:HA3	2.16	0.46
1:A:82:VAL:HG13	1:A:105:VAL:HG13	1.98	0.46
1:A:155:THR:OG1	1:A:310:PHE:CZ	2.64	0.46
1:A:277:ILE:CG2	1:A:284:LEU:HB3	2.46	0.46
1:B:55:LYS:HG3	1:B:115:TYR:OH	2.15	0.46
1:A:89:ASP:OD1	1:A:90:LEU:N	2.40	0.45
1:A:14:ILE:HD11	1:A:275:GLN:CG	2.46	0.45
1:A:5:ASN:N	1:A:5:ASN:HD22	2.14	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:LEU:HD12	1:A:153:LEU:O	2.16	0.45
1:A:221:THR:HB	1:A:300:ILE:CB	2.40	0.45
1:B:247:THR:CG2	1:B:248:LEU:N	2.80	0.45
1:B:21:GLU:HB3	1:B:26:GLN:O	2.16	0.45
1:B:47:CYS:SG	1:B:106:ILE:O	2.75	0.45
1:B:47:CYS:HB2	1:B:106:ILE:HA	1.99	0.45
1:B:101:LYS:HD3	1:B:136:VAL:HG22	1.99	0.45
1:A:103:ILE:N	1:A:103:ILE:HD12	2.32	0.45
1:B:163:LYS:CE	1:B:327:LYS:HG3	2.46	0.45
1:B:165:THR:HG23	1:B:166:GLY:N	2.30	0.45
1:A:15:MET:HA	3:A:925:HOH:O	2.17	0.45
1:A:44:SER:C	1:A:46:LYS:N	2.73	0.45
1:A:49:THR:HG23	1:A:109:ASN:OD1	2.17	0.45
1:B:29:PHE:CZ	1:B:42:VAL:HG21	2.52	0.45
1:B:153:LEU:C	1:B:153:LEU:CD1	2.88	0.45
1:A:277:ILE:O	1:A:280:VAL:HB	2.17	0.44
1:A:29:PHE:HB2	1:A:31:PHE:CE1	2.52	0.44
1:B:217:THR:HG22	1:B:219:ALA:H	1.83	0.44
1:A:206:LEU:HD21	1:A:297:PRO:HB3	2.00	0.44
1:A:218:SER:HA	1:A:303:ASP:OD2	2.17	0.44
1:B:269:GLU:CD	1:B:269:GLU:H	2.25	0.44
1:A:204:ILE:HD13	1:A:229:LYS:HB2	2.00	0.44
1:A:217:THR:HA	2:A:347:IVS:O36	2.18	0.44
1:B:23:GLY:HA2	1:B:89:ASP:CG	2.43	0.44
1:B:142:GLU:O	1:B:146:GLN:CG	2.59	0.44
1:B:222:VAL:CG2	1:B:226:PHE:CD1	3.00	0.44
1:B:137:ASP:HB3	3:B:833:HOH:O	2.16	0.44
1:B:246:VAL:HG22	1:B:286:MET:HE2	1.98	0.44
1:B:223:PRO:HG2	1:B:226:PHE:CB	2.46	0.44
1:A:210:ASN:O	1:A:298:THR:HG23	2.18	0.44
1:B:229:LYS:O	1:B:232:GLN:HB2	2.18	0.44
1:B:318:HIS:CD2	1:B:318:HIS:N	2.86	0.44
1:B:327:LYS:O	1:B:327:LYS:CG	2.65	0.43
1:A:65:THR:HG21	1:A:90:LEU:HB2	2.00	0.43
1:A:129:LYS:HB2	1:A:138:PRO:HD3	2.00	0.43
1:B:89:ASP:OD1	1:B:90:LEU:N	2.49	0.43
1:B:125:GLY:O	1:B:126:LEU:HG	2.17	0.43
1:A:294:PHE:HB3	1:A:295:PRO:CD	2.48	0.43
1:B:58:TYR:HD2	1:B:104:GLU:HG3	1.82	0.43
1:A:159:PRO:HD3	1:A:166:GLY:HA3	2.00	0.43
1:B:146:GLN:HB3	1:B:148:LYS:HZ3	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:THR:HG22	1:A:51:GLY:H	1.83	0.43
1:B:58:TYR:CZ	1:B:60:SER:HA	2.53	0.43
1:B:155:THR:HG23	1:B:156:PHE:N	2.33	0.43
1:B:277:ILE:HD13	1:B:277:ILE:HA	1.70	0.43
1:A:38:ALA:HB3	1:A:132:SER:N	2.33	0.43
1:B:21:GLU:CD	1:B:21:GLU:H	2.27	0.43
1:A:201:VAL:CG2	1:A:226:PHE:CE1	3.01	0.43
1:B:194:GLN:HA	1:B:211:CYS:O	2.19	0.43
1:B:222:VAL:CG2	1:B:223:PRO:HD2	2.49	0.42
1:B:272:TYR:CE1	1:B:308:LYS:HB2	2.54	0.42
1:A:82:VAL:HG13	1:A:105:VAL:HG11	1.99	0.42
1:A:155:THR:HG23	1:A:310:PHE:HE1	1.83	0.42
1:B:221:THR:HG23	1:B:292:LEU:HB3	2.01	0.42
1:A:234:LEU:CA	3:A:924:HOH:O	2.67	0.42
1:A:242:LEU:C	1:A:244:PHE:H	2.27	0.42
1:B:35:THR:HA	1:B:126:LEU:HB2	2.00	0.42
1:A:14:ILE:CG2	1:A:218:SER:HB2	2.50	0.42
1:A:44:SER:HB2	1:A:104:GLU:HG2	2.01	0.42
1:B:294:PHE:HB3	1:B:295:PRO:HD2	2.02	0.42
1:B:25:ASN:O	1:B:26:GLN:CB	2.62	0.42
1:B:296:VAL:O	1:B:297:PRO:C	2.60	0.42
1:A:119:THR:O	1:A:119:THR:CG2	2.68	0.42
1:A:294:PHE:HB3	1:A:295:PRO:HD2	2.02	0.42
1:B:24:ASP:OD2	1:B:64:ARG:HB2	2.20	0.42
1:B:68:LYS:HE3	1:B:68:LYS:HB2	1.89	0.42
1:B:134:GLY:O	1:B:135:SER:C	2.63	0.42
1:A:21:GLU:N	1:A:21:GLU:CD	2.78	0.42
1:A:40:LEU:O	1:A:102:PHE:HB2	2.20	0.42
1:A:155:THR:HG23	1:A:310:PHE:CE1	2.55	0.42
1:A:215:SER:HB3	1:A:306:MET:HE1	2.02	0.41
1:A:274:LEU:HD23	1:A:287:LEU:HD22	2.02	0.41
1:B:269:GLU:CD	1:B:269:GLU:N	2.78	0.41
1:B:270:PRO:HA	1:B:273:TYR:CE2	2.54	0.41
1:B:149:ILE:HD11	1:B:170:ILE:O	2.21	0.41
1:B:222:VAL:O	1:B:291:GLY:HA2	2.20	0.41
1:B:307:ARG:HD2	3:B:912:HOH:O	2.20	0.41
1:A:11:PHE:CZ	1:A:12:GLN:NE2	2.88	0.41
1:B:6:ILE:HB	1:B:168:LEU:HB3	2.01	0.41
1:B:66:TYR:O	1:B:67:GLU:HB2	2.18	0.41
1:A:33:LEU:HD13	1:A:168:LEU:HD22	2.02	0.41
1:A:65:THR:O	1:A:67:GLU:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:ASN:O	1:A:251:ASN:C	2.63	0.41
1:B:257:PHE:HB3	1:B:259:PHE:CE2	2.56	0.41
1:A:78:VAL:N	3:A:927:HOH:O	2.26	0.41
1:B:290:ILE:HG13	2:B:348:IVS:HC61	2.01	0.41
1:B:84:GLY:HA3	1:B:104:GLU:O	2.21	0.41
1:B:251:ASN:CB	1:B:254:LEU:HG	2.47	0.41
1:A:26:GLN:HG2	1:A:28:PRO:HD3	2.03	0.41
1:A:226:PHE:O	1:A:230:MET:HB2	2.20	0.41
1:A:238:LYS:HG3	1:A:245:TYR:CE1	2.56	0.41
1:B:182:LEU:HD21	1:B:312:VAL:HG21	2.03	0.41
1:A:112:GLU:HB3	1:A:113:PRO:HA	2.03	0.41
1:A:188:ASN:C	1:A:188:ASN:ND2	2.75	0.41
1:B:176:ARG:O	1:B:326:LYS:HD2	2.20	0.41
1:B:182:LEU:CD1	1:B:323:ALA:HB2	2.51	0.40
1:A:326:LYS:C	1:A:328:ASN:N	2.78	0.40
1:B:220:ILE:HD13	1:B:220:ILE:HA	1.97	0.40
1:B:222:VAL:HG22	1:B:223:PRO:HD2	2.04	0.40
1:A:198:ASP:O	1:A:259:PHE:HA	2.21	0.40
1:A:290:ILE:HG13	2:A:347:IVS:HC61	2.03	0.40
1:B:60:SER:HB2	1:B:66:TYR:CD1	2.56	0.40
1:B:88:LYS:HZ1	1:B:99:PRO:HB2	1.86	0.40
1:B:163:LYS:HB3	1:B:327:LYS:HE3	2.04	0.40
1:B:191:LEU:HB3	1:B:192:TYR:H	1.74	0.40
1:A:4:ASP:OD2	1:A:94:GLY:HA3	2.21	0.40
1:B:15:MET:HG3	1:B:16:PHE:N	2.36	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	327/329 (99%)	298 (91%)	23 (7%)	6 (2%)	6 18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	327/329 (99%)	293 (90%)	28 (9%)	6 (2%)	6	18
All	All	654/658 (99%)	591 (90%)	51 (8%)	12 (2%)	6	18

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	67	GLU
1	A	26	GLN
1	A	68	LYS
1	A	135	SER
1	B	49	THR
1	B	70	GLY
1	B	135	SER
1	B	161	HIS
1	A	161	HIS
1	A	232	GLN
1	A	297	PRO
1	B	240	PRO

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	294/294 (100%)	264 (90%)	30 (10%)	7	18
1	B	294/294 (100%)	274 (93%)	20 (7%)	14	35
All	All	588/588 (100%)	538 (92%)	50 (8%)	10	25

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	19	ASP
1	A	21	GLU
1	A	40	LEU

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Mol	Chain	Res	Type
1	A	42	VAL
1	A	45	VAL
1	A	54	THR
1	A	62	LYS
1	A	71	THR
1	A	72	LYS
1	A	73	VAL
1	A	93	VAL
1	A	119	THR
1	A	144	LYS
1	A	151	ASN
1	A	153	LEU
1	A	155	THR
1	A	165	THR
1	A	188	ASN
1	A	221	THR
1	A	247	THR
1	A	250	ASN
1	A	251	ASN
1	A	269	GLU
1	A	270	PRO
1	A	275	GLN
1	A	277	ILE
1	A	284	LEU
1	A	287	LEU
1	A	290	ILE
1	B	3	ASN
1	B	21	GLU
1	B	26	GLN
1	B	32	ILE
1	B	73	VAL
1	B	119	THR
1	B	124	LEU
1	B	142	GLU
1	B	149	ILE
1	B	153	LEU
1	B	155	THR
1	B	165	THR
1	B	187	LEU
1	B	221	THR
1	B	233	ASN
1	B	240	PRO

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Mol	Chain	Res	Type
1	B	277	ILE
1	B	282	PRO
1	B	284	LEU
1	B	326	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	26	GLN
1	A	56	HIS
1	A	76	ASN
1	A	145	ASN
1	A	147	ASN
1	A	151	ASN
1	A	161	HIS
1	A	188	ASN
1	A	189	HIS
1	A	200	HIS
1	A	233	ASN
1	A	251	ASN
1	B	203	ASN
1	B	275	GLN
1	B	276	HIS
1	B	288	ASN
1	B	328	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 oligosaccharides in this entry.

5.6 Ligand geometry

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	IVS	B	348	-	33,33,33	1.18	5 (15%)	43,44,44	1.57	6 (13%)
2	IVS	A	347	-	33,33,33	1.03	1 (3%)	43,44,44	1.57	6 (13%)

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	IVS	B	348	-	-	4/47/47/47	-
2	IVS	A	347	-	-	5/47/47/47	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	348	IVS	C52-C53	-2.91	1.45	1.50
2	A	347	IVS	C51-C23	2.89	1.58	1.53
2	B	348	IVS	C51-C23	2.85	1.58	1.53
2	B	348	IVS	C52-C51	2.58	1.56	1.53
2	B	348	IVS	C21-N22	-2.55	1.28	1.34
2	B	348	IVS	O72-C51	2.37	1.48	1.43

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	347	IVS	O56-C53-C52	6.67	123.46	111.17
2	B	348	IVS	O56-C53-C52	6.30	122.78	111.17
2	B	348	IVS	O75-C53-C52	-3.64	116.20	124.65
2	A	347	IVS	O75-C53-C52	-3.59	116.31	124.65
2	A	347	IVS	C18-C17-N16	-2.94	102.34	110.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	348	IVS	C18-C17-N16	-2.51	103.51	110.32
2	B	348	IVS	C21-C20-N19	-2.44	103.69	110.32
2	A	347	IVS	C17-N16-C12	-2.43	116.99	121.80
2	B	348	IVS	C10-C12-N16	-2.32	113.08	116.25
2	B	348	IVS	C23-N22-C21	-2.23	119.38	123.25
2	A	347	IVS	C10-C12-N16	-2.20	113.25	116.25
2	A	347	IVS	O56-C53-O75	-2.15	118.25	123.63

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	347	IVS	O72-C51-C52-C53
2	A	347	IVS	C23-C51-C52-C53
2	A	347	IVS	C51-C23-C58-C60
2	B	348	IVS	C51-C23-C58-C60
2	A	347	IVS	N22-C23-C58-C60
2	B	348	IVS	N22-C23-C58-C60
2	B	348	IVS	C52-C53-O56-C57
2	A	347	IVS	C52-C53-O56-C57
2	B	348	IVS	C51-C52-C53-O56

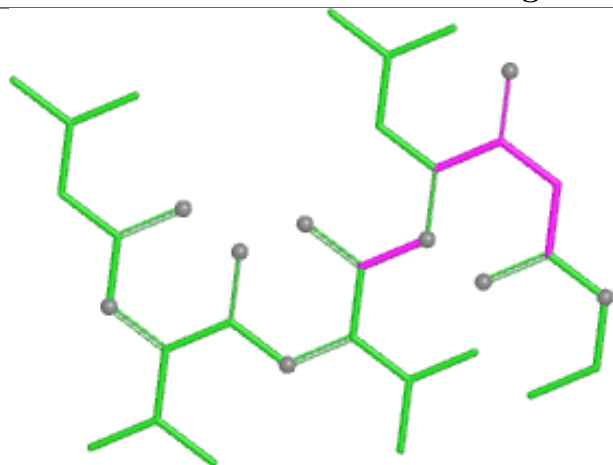
There are no ring outliers.

2 monomers are involved in 3 short contacts:

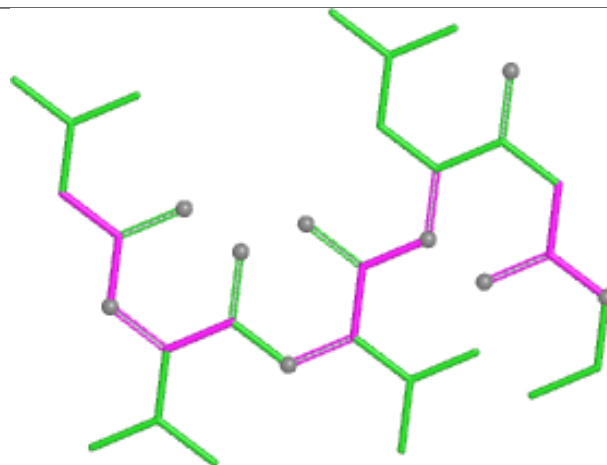
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	348	IVS	1	0
2	A	347	IVS	2	0

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.

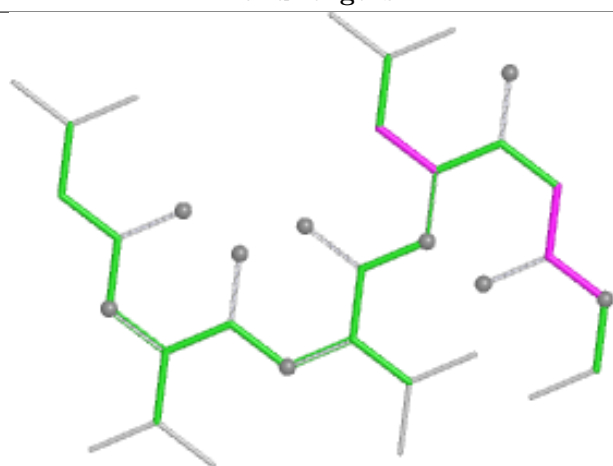
Ligand IVS B 348



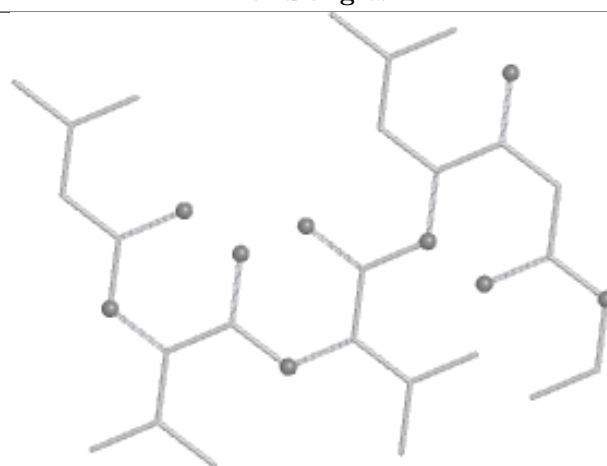
Bond lengths



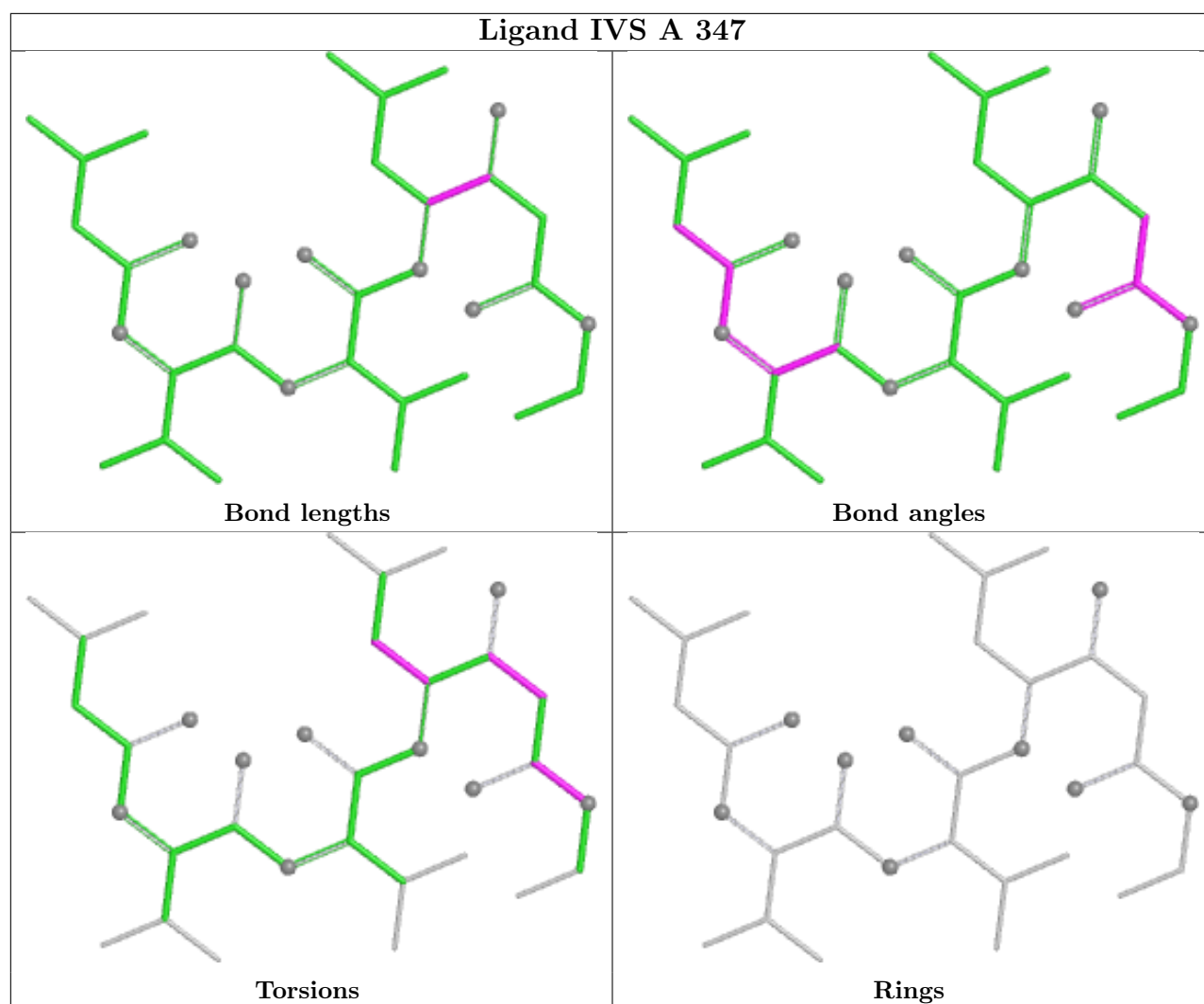
Bond angles



Torsions



Rings



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.