



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 29, 2026 – 02:20 PM JST

PDB ID : 9WNT / pdb_00009wnt
Title : Crystal structure of SN45 TCR in complex with C14-Gly-Gly-Ser-Ile lipopeptide-bound Mamu-B*05104
Authors : Morita, D.; Sugita, M.
Deposited on : 2025-09-05
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 : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

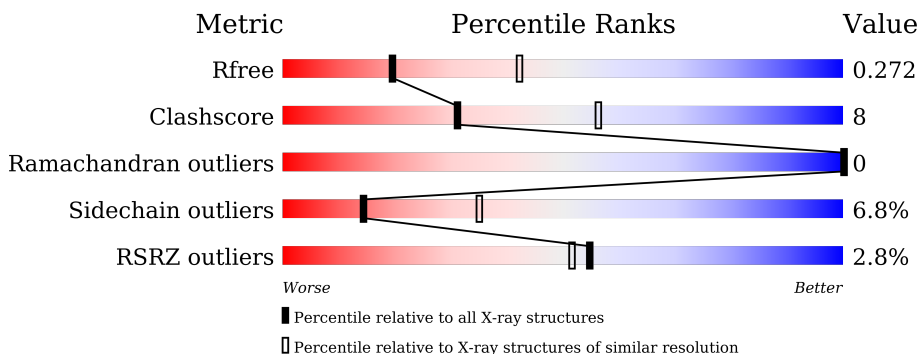
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

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



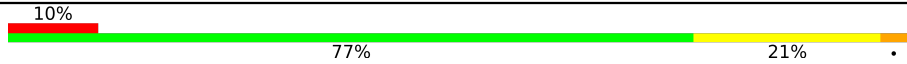



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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

Mol	Chain	Length	Quality of chain
1	A	276	
1	F	276	
2	B	100	
2	G	100	
3	C	4	
3	H	4	

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Mol	Chain	Length	Quality of chain
4	D	197	 10% 77% 21% .
4	I	197	 79% 18% .
5	E	245	 4% 72% 26% .
5	J	245	 74% 24% .

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 13307 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 MHC class I antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	276	Total	C	N	O	S	0	1	0
			2260	1411	412	431	6			
1	F	276	Total	C	N	O	S	0	1	0
			2260	1411	412	431	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	128	GLU	ARG	conflict	UNP B2ZHY7
A	177	GLU	LYS	conflict	UNP B2ZHY7
A	223	GLU	ASP	conflict	UNP B2ZHY7
A	264	GLU	LYS	conflict	UNP B2ZHY7
F	128	GLU	ARG	conflict	UNP B2ZHY7
F	177	GLU	LYS	conflict	UNP B2ZHY7
F	223	GLU	ASP	conflict	UNP B2ZHY7
F	264	GLU	LYS	conflict	UNP B2ZHY7

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	100	Total	C	N	O	S	0	0	0
			827	526	141	156	4			
2	G	100	Total	C	N	O	S	0	0	0
			827	526	141	156	4			

- Molecule 3 is a protein called 4-mer Lipopeptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	4	Total	C	N	O	0	0	0
			23	13	4	6			
3	H	4	Total	C	N	O	0	0	0
			23	13	4	6			

- Molecule 4 is a protein called SN45 T cell receptor alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	197	Total 1495	C 946	N 247	O 293	S 9	0	0	0
4	I	197	Total 1526	C 962	N 252	O 302	S 10	0	0	0

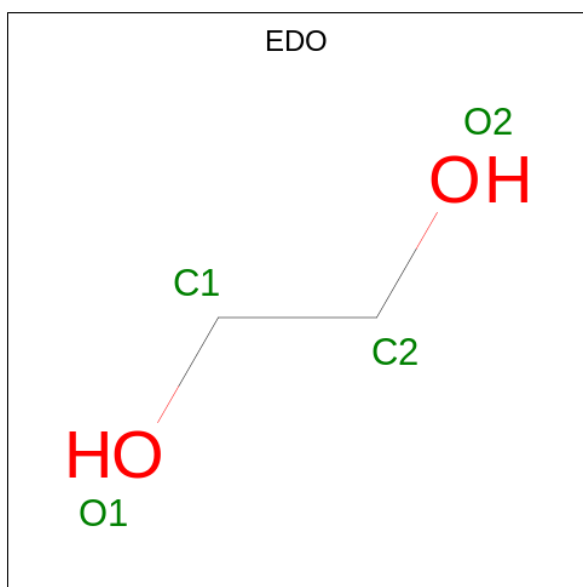
- Molecule 5 is a protein called SN45 T cell receptor beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	E	245	Total 1949	C 1244	N 338	O 359	S 8	0	0	0
5	J	245	Total 1980	C 1259	N 338	O 375	S 8	0	0	0

- Molecule 6 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	2	Total I 2 2	0	0
6	B	2	Total I 2 2	0	0
6	D	1	Total I 1 1	0	0
6	E	2	Total I 2 2	0	0
6	F	2	Total I 2 2	0	0
6	G	1	Total I 1 1	0	0
6	I	3	Total I 3 3	0	0
6	J	5	Total I 5 5	0	0

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).

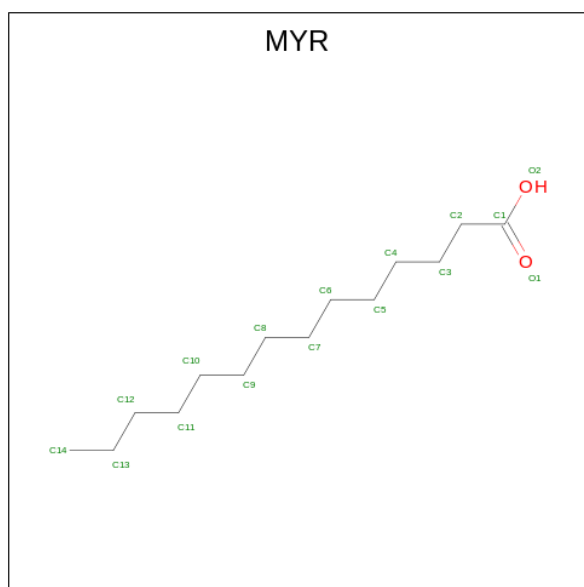


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 4 2 2	0	0
7	A	1	Total C O 4 2 2	0	0
7	A	1	Total C O 4 2 2	0	0
7	A	1	Total C O 4 2 2	0	0
7	B	1	Total C O 4 2 2	0	0
7	B	1	Total C O 4 2 2	0	0
7	B	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0
7	F	1	Total C O 4 2 2	0	0
7	I	1	Total C O 4 2 2	0	0
7	J	1	Total C O 4 2 2	0	0
7	J	1	Total C O 4 2 2	0	0

- Molecule 8 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	B	1	Total Na 1 1	0	0
8	G	1	Total Na 1 1	0	0
8	I	1	Total Na 1 1	0	0

- Molecule 9 is MYRISTIC ACID (CCD ID: MYR) (formula: C₁₄H₂₈O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	C	1	Total C O 15 14 1	0	0
9	H	1	Total C O 15 14 1	0	0

- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	5	Total O 5 5	0	0
10	B	6	Total O 6 6	0	0
10	E	2	Total O 2 2	0	0
10	F	5	Total O 5 5	0	0

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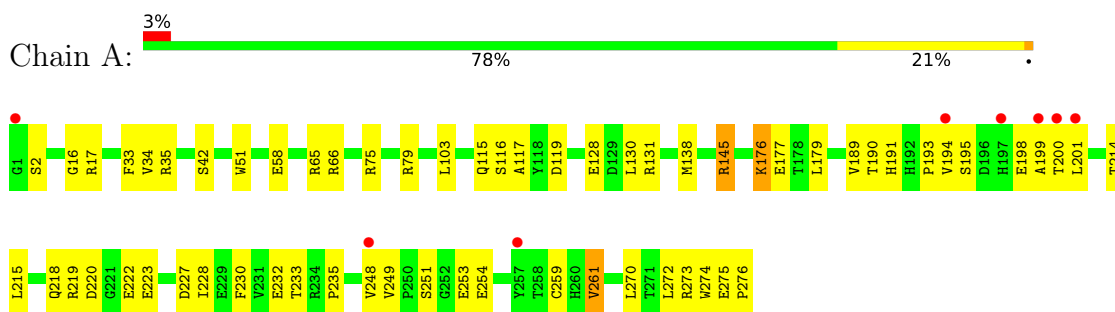
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	G	3	Total O 3 3	0	0
10	I	9	Total O 9 9	0	0
10	J	4	Total O 4 4	0	0

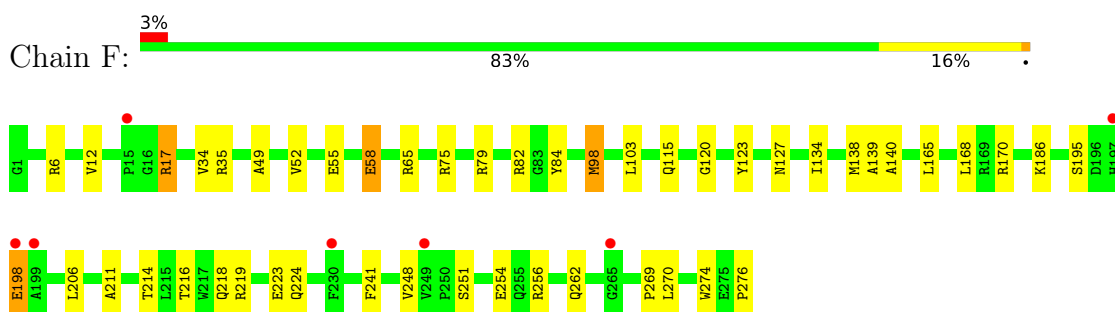
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

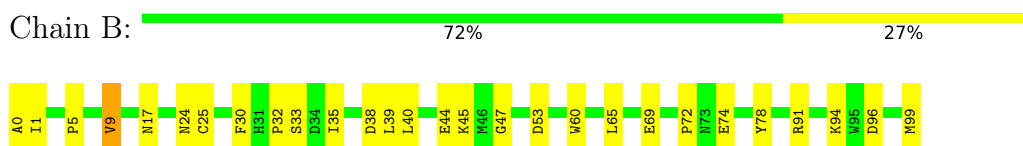
- Molecule 1: MHC class I antigen



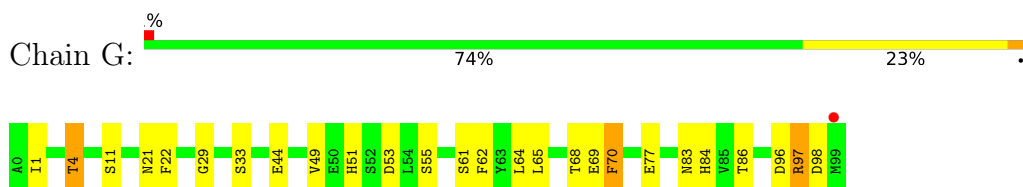
- Molecule 1: MHC class I antigen



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: 4-mer Lipopeptide

Chain C:  50% 50%




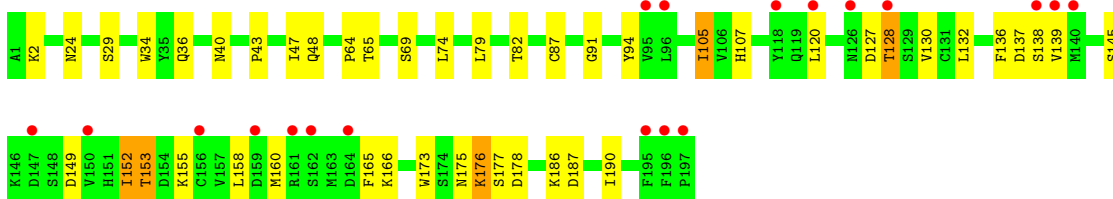
- Molecule 3: 4-mer Lipopeptide


Chain H:  75% 25%




- Molecule 4: SN45 T cell receptor alpha chain

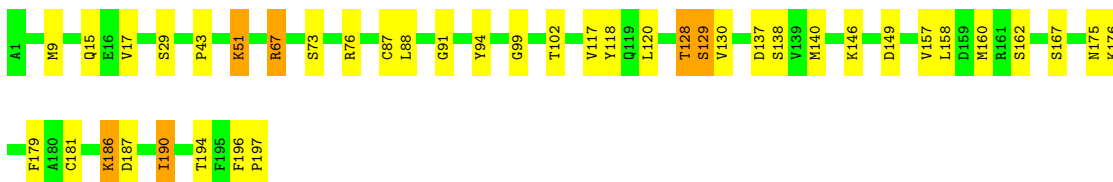
Chain D:  10% 77% 21%







- Molecule 4: SN45 T cell receptor alpha chain

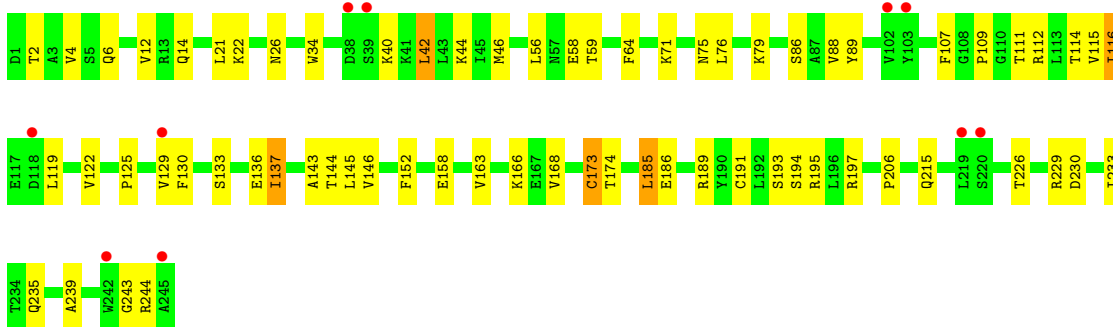
Chain I:  79% 18%




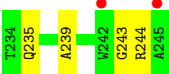


- Molecule 5: SN45 T cell receptor beta chain

Chain E:  4% 72% 26%

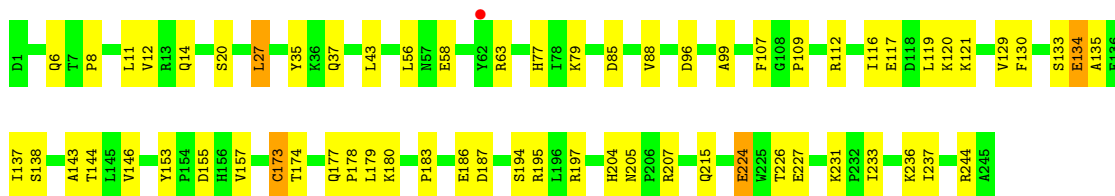






- Molecule 5: SN45 T cell receptor beta chain

Chain J:  74% 24%



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.08Å 94.62Å 123.05Å 90.00° 90.47° 90.00°	Depositor
Resolution (Å)	49.46 – 2.70 49.46 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.1 (49.46-2.70) 99.1 (49.46-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 2.69Å)	Xtrriage
Refinement program	PHENIX 2.0	Depositor
R, R_{free}	0.204 , 0.274 0.206 , 0.272	Depositor DCC
R_{free} test set	2642 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	61.6	Xtrriage
Anisotropy	0.125	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.018 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13307	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, IOD, NA, MYR

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	0.45	0/2327	0.64	0/3164
1	F	0.45	0/2327	0.66	0/3164
2	B	0.43	0/852	0.65	0/1153
2	G	0.41	0/852	0.61	0/1153
3	C	0.91	0/22	1.21	0/26
3	H	0.92	0/22	1.06	0/26
4	D	0.47	0/1531	0.72	1/2087 (0.0%)
4	I	0.50	0/1562	0.75	0/2125
5	E	0.44	0/2004	0.67	0/2721
5	J	0.50	0/2035	0.73	0/2761
All	All	0.46	0/13534	0.69	1/18380 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
4	I	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	139	VAL	N-CA-C	-6.27	106.21	112.17

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	I	67	ARG	Sidechain

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	2260	0	2097	38	0
1	F	2260	0	2097	26	0
2	B	827	0	788	18	0
2	G	827	0	788	16	0
3	C	23	0	21	0	0
3	H	23	0	21	0	0
4	D	1495	0	1401	33	0
4	I	1526	0	1460	24	0
5	E	1949	0	1872	42	0
5	J	1980	0	1914	38	0
6	A	2	0	0	0	0
6	B	2	0	0	1	0
6	D	1	0	0	0	0
6	E	2	0	0	0	0
6	F	2	0	0	0	0
6	G	1	0	0	0	0
6	I	3	0	0	1	0
6	J	5	0	0	0	0
7	A	16	0	24	2	0
7	B	12	0	18	1	0
7	D	8	0	12	3	0
7	F	4	0	6	0	0
7	I	4	0	6	0	0
7	J	8	0	12	2	0
8	B	1	0	0	0	0
8	G	1	0	0	0	0
8	I	1	0	0	0	0
9	C	15	0	27	0	0
9	H	15	0	27	1	0
10	A	5	0	0	1	0
10	B	6	0	0	0	0
10	E	2	0	0	0	0
10	F	5	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	G	3	0	0	0	0
10	I	9	0	0	1	0
10	J	4	0	0	0	0
All	All	13307	0	12591	211	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 8.

All (211) 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)
4:D:137:ASP:H	7:D:202:EDO:H11	1.34	0.93
2:B:17:ASN:HD21	2:B:74:GLU:HG2	1.41	0.85
4:D:132:LEU:HD22	5:E:144:THR:HG21	1.57	0.85
5:E:6:GLN:HB2	5:E:109:PRO:HD2	1.57	0.83
2:G:4:THR:HA	2:G:86:THR:HG21	1.60	0.82
4:D:82:THR:HG23	4:D:105:ILE:HA	1.61	0.81
4:I:190:ILE:HD11	4:I:194:THR:HG21	1.65	0.78
1:A:191:HIS:NE2	1:A:199:ALA:HB1	2.01	0.75
1:A:191:HIS:CD2	1:A:199:ALA:HB1	2.22	0.74
4:D:158:LEU:HD13	5:E:197:ARG:HB2	1.71	0.72
5:J:37:GLN:HB2	5:J:43:LEU:HD23	1.71	0.72
1:A:261:VAL:HG22	1:A:270:LEU:HB3	1.71	0.72
4:I:130:VAL:HG11	5:J:146:VAL:HG21	1.72	0.71
5:E:116:ILE:HD11	5:E:119:LEU:HD13	1.71	0.71
1:A:259:CYS:HB3	1:A:272:LEU:HB2	1.73	0.71
5:J:88:VAL:HG22	5:J:112:ARG:HG3	1.73	0.70
5:E:129:VAL:HG23	5:E:239:ALA:HB3	1.75	0.68
1:F:79:ARG:HD2	1:F:82:ARG:NH2	2.09	0.67
2:G:84:HIS:ND1	2:G:86:THR:HG22	2.09	0.67
1:F:6:ARG:HD3	1:F:98:MET:HE2	1.75	0.67
1:A:194:VAL:HG22	1:A:195:SER:H	1.59	0.67
4:D:145:SER:HB3	4:D:152:ILE:HD11	1.78	0.65
1:A:274:TRP:NE1	1:A:276:PRO:HB3	2.12	0.64
5:E:14:GLN:HB2	5:E:119:LEU:HD21	1.80	0.63
1:F:79:ARG:HH11	1:F:82:ARG:HH22	1.46	0.62
4:D:40:ASN:HA	5:E:112:ARG:HH22	1.64	0.62
2:B:47:GLY:HA2	5:J:186:GLU:HG2	1.82	0.61
1:A:249:VAL:HG21	1:A:254:GLU:HG3	1.80	0.61
1:F:79:ARG:NH1	1:F:82:ARG:HH22	1.98	0.61
4:I:99:GLY:HA3	10:I:302:HOH:O	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:VAL:HG23	1:A:272:LEU:HD12	1.83	0.61
4:D:137:ASP:N	7:D:202:EDO:H11	2.11	0.60
4:I:9:MET:HE3	4:I:102:THR:HG21	1.83	0.60
5:E:206:PRO:HA	5:E:243:GLY:O	2.01	0.60
4:I:29:SER:O	4:I:91:GLY:HA3	2.01	0.60
5:J:231:LYS:HB3	5:J:233:ILE:HG23	1.84	0.60
5:J:205:ASN:OD1	5:J:207:ARG:HG2	2.02	0.60
2:G:21:ASN:HB3	2:G:70:PHE:HE1	1.65	0.59
1:F:35:ARG:HD3	2:G:53:ASP:OD2	2.03	0.58
5:E:56:LEU:HD21	5:E:58:GLU:HG3	1.84	0.58
1:F:58:GLU:CD	1:F:58:GLU:H	2.10	0.58
5:E:64:PHE:HD1	5:E:76:LEU:HD21	1.69	0.58
5:E:233:ILE:O	5:E:235:GLN:HG2	2.03	0.58
5:E:136:GLU:OE1	5:E:144:THR:HG22	2.04	0.58
5:J:35:TYR:HB3	5:J:43:LEU:HD22	1.85	0.58
1:F:103:LEU:HD13	1:F:165:LEU:HD22	1.86	0.58
1:A:119:ASP:HB3	2:B:0:ALA:HA	1.85	0.57
4:D:2:LYS:HE3	4:D:24:ASN:H	1.69	0.57
1:F:75:ARG:O	1:F:79:ARG:HD3	2.05	0.57
1:F:103:LEU:HB2	1:F:168:LEU:HD23	1.86	0.57
2:G:21:ASN:HB3	2:G:70:PHE:CE1	2.40	0.57
2:B:72:PRO:HB2	6:B:102:IOD:I	2.75	0.57
5:E:174:THR:HG23	5:E:194:SER:HB2	1.87	0.56
5:J:144:THR:OG1	5:J:197:ARG:HG3	2.05	0.56
1:F:262:GLN:OE1	1:F:269:PRO:HG3	2.06	0.56
1:A:2:SER:HB2	1:A:103:LEU:O	2.07	0.55
1:A:228:ILE:HD12	1:A:230:PHE:HE2	1.70	0.55
9:H:101:MYR:H21	5:J:99:ALA:HB3	1.87	0.55
5:E:133:SER:HB3	5:E:136:GLU:H	1.71	0.55
4:D:120:LEU:HD11	4:D:132:LEU:HB2	1.88	0.55
5:J:174:THR:HG23	5:J:194:SER:HB2	1.88	0.55
1:A:75:ARG:HG2	1:A:79:ARG:HH21	1.72	0.55
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.42	0.54
1:F:195:SER:HB2	1:F:198:GLU:HB2	1.90	0.54
5:J:20:SER:HB2	5:J:77:HIS:CD2	2.43	0.54
1:F:274:TRP:CZ2	1:F:276:PRO:HB3	2.42	0.54
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.90	0.54
4:I:43:PRO:HD2	5:J:107:PHE:CD2	2.43	0.54
4:D:128:THR:HB	4:D:175:ASN:HB2	1.89	0.53
5:E:163:VAL:O	5:E:166:LYS:HG2	2.09	0.53
5:J:121:LYS:HE2	5:J:187:ASP:HB3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:9:VAL:O	7:B:106:EDO:H21	2.09	0.53
5:E:158:GLU:HB2	5:E:215:GLN:HB3	1.91	0.53
1:F:219:ARG:HB3	1:F:224:GLN:HE22	1.74	0.53
4:D:64:PRO:HD2	4:D:69:SER:O	2.09	0.53
5:E:22:LYS:HE2	5:E:75:ASN:HD21	1.74	0.52
5:E:125:PRO:HB3	5:E:152:PHE:CD1	2.45	0.52
5:E:22:LYS:HE2	5:E:75:ASN:ND2	2.24	0.52
5:E:46:MET:HE1	5:E:89:TYR:HD2	1.74	0.52
5:E:152:PHE:CE1	5:E:191:CYS:HA	2.45	0.51
5:J:137:ILE:HG12	5:J:143:ALA:HB2	1.92	0.51
2:B:17:ASN:ND2	2:B:74:GLU:HG2	2.19	0.51
5:J:204:HIS:HA	5:J:244:ARG:O	2.11	0.51
5:J:6:GLN:HB2	5:J:109:PRO:HD2	1.91	0.51
4:D:2:LYS:HE3	4:D:24:ASN:N	2.25	0.50
5:E:86:SER:HG	5:E:115:VAL:H	1.57	0.50
4:D:79:LEU:HD11	4:D:166:LYS:HD2	1.92	0.50
4:D:29:SER:O	4:D:91:GLY:HA3	2.11	0.50
1:A:191:HIS:CE1	1:A:193:PRO:HD3	2.47	0.50
1:F:120:GLY:HA3	2:G:1:ILE:HG22	1.92	0.50
5:E:40:LYS:HG3	5:E:42:LEU:HB2	1.93	0.50
1:F:84:TYR:HB3	1:F:139:ALA:HB1	1.94	0.50
4:I:157:VAL:O	5:J:173:CYS:SG	2.69	0.50
1:A:115:GLN:HG3	10:A:404:HOH:O	2.11	0.49
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.94	0.49
5:E:146:VAL:HG12	5:E:195:ARG:HG2	1.94	0.49
4:I:138:SER:HB2	6:I:202:IOD:I	2.83	0.49
4:D:47:ILE:HA	7:D:203:EDO:H21	1.94	0.49
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.93	0.49
5:E:2:THR:HA	5:E:26:ASN:HD21	1.78	0.49
4:D:158:LEU:CD2	5:E:173:CYS:HB2	2.43	0.49
4:I:129:SER:HB2	4:I:179:PHE:HD2	1.78	0.49
5:J:130:PHE:HB2	5:J:146:VAL:HG22	1.94	0.48
4:D:107:HIS:HB3	4:D:138:SER:CB	2.43	0.48
5:J:183:PRO:HB3	7:J:307:EDO:H11	1.95	0.48
5:E:12:VAL:HA	5:E:114:THR:O	2.13	0.48
2:G:49:VAL:HG22	2:G:68:THR:HB	1.94	0.48
1:F:55:GLU:CD	1:F:170:ARG:HH12	2.21	0.48
5:J:195:ARG:HD3	7:J:306:EDO:O2	2.13	0.48
5:E:88:VAL:HG22	5:E:112:ARG:HG2	1.96	0.47
4:I:130:VAL:HG11	5:J:146:VAL:CG2	2.44	0.47
4:I:149:ASP:HB2	4:I:176:LYS:HD2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:GLN:HE21	1:A:223:GLU:HG2	1.79	0.47
4:D:176:LYS:HG2	4:D:177:SER:H	1.78	0.47
4:I:118:TYR:CD1	4:I:197:PRO:HG3	2.49	0.47
1:A:66:ARG:HH21	7:A:303:EDO:H12	1.80	0.47
5:E:4:VAL:O	5:E:109:PRO:HD3	2.14	0.47
4:I:43:PRO:HD2	5:J:107:PHE:CG	2.49	0.47
4:D:186:LYS:HG2	4:D:187:ASP:H	1.80	0.47
1:A:176:LYS:HG3	1:A:177:GLU:N	2.30	0.47
5:J:236:LYS:O	5:J:237:ILE:HD13	2.15	0.46
4:D:158:LEU:HD23	4:D:158:LEU:HA	1.78	0.46
5:E:145:LEU:O	5:E:195:ARG:HA	2.15	0.46
4:I:196:PHE:O	4:I:197:PRO:C	2.57	0.46
1:F:214:THR:OG1	1:F:262:GLN:HB3	2.16	0.46
4:I:137:ASP:OD2	4:I:140:MET:HG2	2.16	0.46
4:I:146:LYS:HG2	4:I:187:ASP:OD1	2.15	0.46
1:F:65:ARG:HD2	4:I:94:TYR:CZ	2.51	0.46
1:A:65:ARG:HD2	4:D:94:TYR:CZ	2.50	0.46
1:A:219:ARG:O	1:A:220:ASP:HB2	2.16	0.46
5:E:137:ILE:HD11	5:E:143:ALA:HB2	1.97	0.46
4:D:130:VAL:HG12	4:D:173:TRP:HB3	1.98	0.46
4:D:186:LYS:O	4:D:187:ASP:HB2	2.16	0.45
1:A:218:GLN:HG2	1:A:223:GLU:HG2	1.98	0.45
5:E:34:TRP:HB3	5:E:46:MET:HE3	1.97	0.45
5:J:14:GLN:HG3	5:J:117:GLU:O	2.16	0.45
1:A:145:ARG:CZ	1:A:145:ARG:HB3	2.47	0.45
1:A:201:LEU:HD11	1:A:254:GLU:HB3	1.98	0.45
2:G:11:SER:HA	2:G:22:PHE:O	2.17	0.45
5:E:206:PRO:HD3	5:E:244:ARG:HA	1.97	0.45
2:B:38:ASP:OD1	2:B:45:LYS:HE3	2.16	0.44
4:I:186:LYS:HA	4:I:186:LYS:HD2	1.58	0.44
1:A:191:HIS:NE2	1:A:199:ALA:CB	2.76	0.44
1:A:198:GLU:HA	1:A:251:SER:HB2	1.99	0.44
1:A:232:GLU:HG2	1:A:233:THR:O	2.17	0.44
1:F:127:ASN:OD1	1:F:134:ILE:HD11	2.16	0.44
2:G:33:SER:HB3	2:G:62:PHE:CE2	2.51	0.44
4:D:136:PHE:HE2	4:D:155:LYS:HD3	1.82	0.44
1:F:251:SER:HA	1:F:254:GLU:CD	2.43	0.44
4:I:128:THR:HB	4:I:175:ASN:OD1	2.18	0.44
4:D:120:LEU:HD21	5:E:144:THR:HG23	2.00	0.44
1:F:262:GLN:NE2	10:F:401:HOH:O	2.51	0.44
4:I:117:VAL:HG12	4:I:181:CYS:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:GLU:O	1:A:130:LEU:HG	2.18	0.43
2:G:97:ARG:CZ	2:G:97:ARG:H	2.31	0.43
1:A:273:ARG:O	1:A:274:TRP:C	2.61	0.43
1:A:189:VAL:CG2	1:A:272:LEU:HB3	2.49	0.43
1:A:218:GLN:HA	1:A:222:GLU:O	2.19	0.43
5:J:134:GLU:HG3	5:J:135:ALA:N	2.34	0.43
5:J:224:GLU:HG3	5:J:226:THR:HG23	2.01	0.43
2:G:4:THR:HA	2:G:86:THR:CG2	2.39	0.43
1:A:235:PRO:HG2	2:B:65:LEU:HD13	2.00	0.42
5:J:116:ILE:HG21	5:J:153:TYR:HE2	1.84	0.42
5:J:120:LYS:HD3	5:J:227:GLU:OE2	2.19	0.42
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.55	0.42
4:D:40:ASN:HA	5:E:112:ARG:NH2	2.33	0.42
4:D:160:MET:HG3	4:D:165:PHE:HB3	2.01	0.42
5:E:40:LYS:O	5:E:42:LEU:HD12	2.19	0.42
2:B:91:ARG:HE	2:B:91:ARG:HB2	1.71	0.42
1:F:218:GLN:HG2	1:F:223:GLU:HG2	2.00	0.42
5:E:21:LEU:HD22	5:E:111:THR:HG21	2.02	0.42
1:A:51:TRP:CZ2	1:A:179:LEU:HD11	2.55	0.42
4:D:153:THR:HG21	5:E:193:SER:OG	2.20	0.42
2:G:96:ASP:C	2:G:98:ASP:H	2.28	0.42
5:J:20:SER:HB2	5:J:77:HIS:NE2	2.35	0.42
5:J:56:LEU:HD21	5:J:58:GLU:HB2	2.02	0.42
1:F:49:ALA:O	1:F:52:VAL:HG12	2.19	0.42
1:F:211:ALA:HB2	1:F:241:PHE:CE2	2.55	0.42
1:A:16:GLY:HA3	7:A:304:EDO:H22	2.02	0.42
1:A:35:ARG:HD3	2:B:53:ASP:OD2	2.20	0.42
4:D:149:ASP:HB3	4:D:176:LYS:HD3	2.01	0.42
2:B:96:ASP:HB3	2:B:99:MET:HB2	2.01	0.41
5:E:229:ARG:HG2	5:E:230:ASP:N	2.35	0.41
4:I:158:LEU:HD11	5:J:197:ARG:HB2	2.03	0.41
4:I:87:CYS:C	4:I:88:LEU:HD12	2.45	0.41
5:J:79:LYS:HE2	5:J:79:LYS:HB3	1.88	0.41
5:E:185:LEU:HG	2:G:44:GLU:HB2	2.03	0.41
4:D:43:PRO:HG2	5:E:107:PHE:CD1	2.55	0.41
4:D:127:ASP:HB2	4:D:175:ASN:HD21	1.86	0.41
4:D:190:ILE:H	4:D:190:ILE:HG13	1.60	0.41
1:F:17:ARG:HD3	1:F:17:ARG:HA	1.92	0.41
1:A:79:ARG:HE	1:A:79:ARG:HB2	1.62	0.41
5:E:130:PHE:HB2	5:E:146:VAL:HG22	2.03	0.41
2:G:51:HIS:HA	2:G:65:LEU:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:J:27:LEU:HD13	5:J:27:LEU:HA	1.82	0.41
1:A:219:ARG:O	1:A:222:GLU:HB3	2.20	0.41
2:B:40:LEU:O	2:B:78:TYR:HA	2.21	0.41
2:G:29:GLY:HA2	2:G:61:SER:OG	2.20	0.41
2:G:64:LEU:HD23	2:G:64:LEU:HA	1.90	0.41
4:I:158:LEU:HD21	4:I:160:MET:HG3	2.03	0.41
4:I:51:LYS:HG2	4:I:67:ARG:HD2	2.02	0.41
5:J:8:PRO:HG2	5:J:11:LEU:HG	2.02	0.41
1:F:123:TYR:CZ	1:F:140:ALA:HA	2.56	0.40
5:J:155:ASP:OD1	5:J:155:ASP:N	2.53	0.40
1:A:251:SER:C	1:A:253:GLU:H	2.28	0.40
5:J:63:ARG:NH2	5:J:85:ASP:OD2	2.54	0.40
5:J:178:PRO:O	5:J:179:LEU:HD23	2.22	0.40
4:D:34:TRP:CZ2	4:D:87:CYS:HB2	2.56	0.40
5:J:157:VAL:HA	5:J:215:GLN:O	2.21	0.40
2:B:5:PRO:CB	2:B:30:PHE:HB3	2.51	0.40
2:B:40:LEU:HA	2:B:44:GLU:O	2.22	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	275/276 (100%)	256 (93%)	19 (7%)	0	100	100
1	F	275/276 (100%)	264 (96%)	11 (4%)	0	100	100
2	B	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
2	G	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
3	C	2/4 (50%)	2 (100%)	0	0	100	100
3	H	2/4 (50%)	2 (100%)	0	0	100	100
4	D	195/197 (99%)	170 (87%)	25 (13%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	I	195/197 (99%)	183 (94%)	12 (6%)	0	100	100
5	E	243/245 (99%)	226 (93%)	17 (7%)	0	100	100
5	J	243/245 (99%)	231 (95%)	12 (5%)	0	100	100
All	All	1626/1644 (99%)	1522 (94%)	104 (6%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	229/229 (100%)	213 (93%)	16 (7%)	14	34
1	F	229/229 (100%)	215 (94%)	14 (6%)	17	40
2	B	93/93 (100%)	86 (92%)	7 (8%)	12	31
2	G	93/93 (100%)	86 (92%)	7 (8%)	12	31
3	C	2/2 (100%)	0	2 (100%)	0	0
3	H	2/2 (100%)	1 (50%)	1 (50%)	0	0
4	D	162/173 (94%)	152 (94%)	10 (6%)	16	39
4	I	173/173 (100%)	161 (93%)	12 (7%)	14	34
5	E	207/218 (95%)	193 (93%)	14 (7%)	14	35
5	J	218/218 (100%)	206 (94%)	12 (6%)	19	45
All	All	1408/1430 (98%)	1313 (93%)	95 (7%)	14	36

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
1	A	42	SER
1	A	58	GLU
1	A	116	SER
1	A	131	ARG

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Mol	Chain	Res	Type
1	A	138	MET
1	A	145	ARG
1	A	176	LYS
1	A	190	THR
1	A	200	THR
1	A	214	THR
1	A	215	LEU
1	A	227	ASP
1	A	248	VAL
1	A	261	VAL
1	A	275	GLU
2	B	1	ILE
2	B	9	VAL
2	B	32	PRO
2	B	33	SER
2	B	35	ILE
2	B	69	GLU
2	B	94	LYS
3	C	4	SER
3	C	5	ILE
4	D	36	GLN
4	D	48	GLN
4	D	65	THR
4	D	74	LEU
4	D	105	ILE
4	D	128	THR
4	D	152	ILE
4	D	153	THR
4	D	176	LYS
4	D	178	ASP
5	E	42	LEU
5	E	44	LYS
5	E	59	THR
5	E	71	LYS
5	E	79	LYS
5	E	116	ILE
5	E	122	VAL
5	E	137	ILE
5	E	168	VAL
5	E	173	CYS
5	E	185	LEU
5	E	186	GLU

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Mol	Chain	Res	Type
5	E	189	ARG
5	E	226	THR
1	F	12	VAL
1	F	17	ARG
1	F	34	VAL
1	F	58	GLU
1	F	98	MET
1	F	115	GLN
1	F	138	MET
1	F	186	LYS
1	F	198	GLU
1	F	206	LEU
1	F	216	THR
1	F	248	VAL
1	F	256	ARG
1	F	270	LEU
2	G	4	THR
2	G	55	SER
2	G	69	GLU
2	G	70	PHE
2	G	77	GLU
2	G	83	ASN
2	G	97	ARG
3	H	5	ILE
4	I	15	GLN
4	I	17	VAL
4	I	51	LYS
4	I	73	SER
4	I	76	ARG
4	I	120	LEU
4	I	128	THR
4	I	129	SER
4	I	162	SER
4	I	167	SER
4	I	186	LYS
4	I	190	ILE
5	J	12	VAL
5	J	27	LEU
5	J	96	ASP
5	J	119	LEU
5	J	129	VAL
5	J	133	SER

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Mol	Chain	Res	Type
5	J	134	GLU
5	J	138	SER
5	J	173	CYS
5	J	177	GLN
5	J	180	LYS
5	J	224	GLU

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	87	GLN
1	A	192	HIS
2	B	13	HIS
2	B	17	ASN
2	B	83	ASN
4	D	15	GLN
4	D	141	ASN
1	F	181	HIS
1	F	260	HIS
2	G	2	GLN
2	G	13	HIS
2	G	73	ASN
4	I	15	GLN
4	I	151	HIS
5	J	14	GLN
5	J	29	HIS
5	J	95	GLN
5	J	105	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 36 ligands modelled in this entry, 21 are monoatomic - leaving 15 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
7	EDO	B	105	-	3,3,3	0.24	0	2,2,2	0.44	0
7	EDO	I	205	-	3,3,3	0.24	0	2,2,2	0.21	0
7	EDO	J	307	-	3,3,3	0.30	0	2,2,2	1.16	0
9	MYR	C	101	3	14,14,15	0.47	0	13,13,15	0.51	0
7	EDO	D	202	-	3,3,3	0.21	0	2,2,2	0.75	0
7	EDO	A	306	-	3,3,3	0.28	0	2,2,2	0.43	0
7	EDO	D	203	-	3,3,3	0.25	0	2,2,2	0.48	0
7	EDO	A	304	-	3,3,3	0.25	0	2,2,2	0.57	0
7	EDO	A	303	-	3,3,3	0.33	0	2,2,2	0.33	0
9	MYR	H	101	3	14,14,15	0.43	0	13,13,15	0.51	0
7	EDO	F	303	-	3,3,3	0.28	0	2,2,2	0.57	0
7	EDO	B	104	-	3,3,3	0.32	0	2,2,2	0.29	0
7	EDO	B	106	-	3,3,3	0.30	0	2,2,2	0.26	0
7	EDO	A	305	-	3,3,3	0.30	0	2,2,2	0.18	0
7	EDO	J	306	8	3,3,3	0.29	0	2,2,2	0.26	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
7	EDO	B	105	-	-	0/1/1/1	-
7	EDO	I	205	-	-	0/1/1/1	-
7	EDO	J	307	-	-	1/1/1/1	-
9	MYR	C	101	3	-	6/11/12/13	-
7	EDO	D	202	-	-	1/1/1/1	-
7	EDO	A	306	-	-	1/1/1/1	-
7	EDO	D	203	-	-	1/1/1/1	-
7	EDO	A	304	-	-	1/1/1/1	-
7	EDO	A	303	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	MYR	H	101	3	-	5/11/12/13	-
7	EDO	F	303	-	-	0/1/1/1	-
7	EDO	B	104	-	-	1/1/1/1	-
7	EDO	B	106	-	-	1/1/1/1	-
7	EDO	A	305	-	-	0/1/1/1	-
7	EDO	J	306	8	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	C	101	MYR	C1-C2-C3-C4
9	C	101	MYR	C10-C11-C12-C13
9	C	101	MYR	C6-C7-C8-C9
9	H	101	MYR	C5-C6-C7-C8
9	C	101	MYR	C4-C5-C6-C7
7	A	304	EDO	O1-C1-C2-O2
7	A	306	EDO	O1-C1-C2-O2
7	D	203	EDO	O1-C1-C2-O2
9	H	101	MYR	C7-C8-C9-C10
9	H	101	MYR	C11-C10-C9-C8
9	C	101	MYR	C11-C10-C9-C8
9	H	101	MYR	C10-C11-C12-C13
9	C	101	MYR	C5-C6-C7-C8
7	B	106	EDO	O1-C1-C2-O2
9	H	101	MYR	C1-C2-C3-C4
7	J	307	EDO	O1-C1-C2-O2
7	B	104	EDO	O1-C1-C2-O2
7	D	202	EDO	O1-C1-C2-O2

There are no ring outliers.

8 monomers are involved in 9 short contacts:

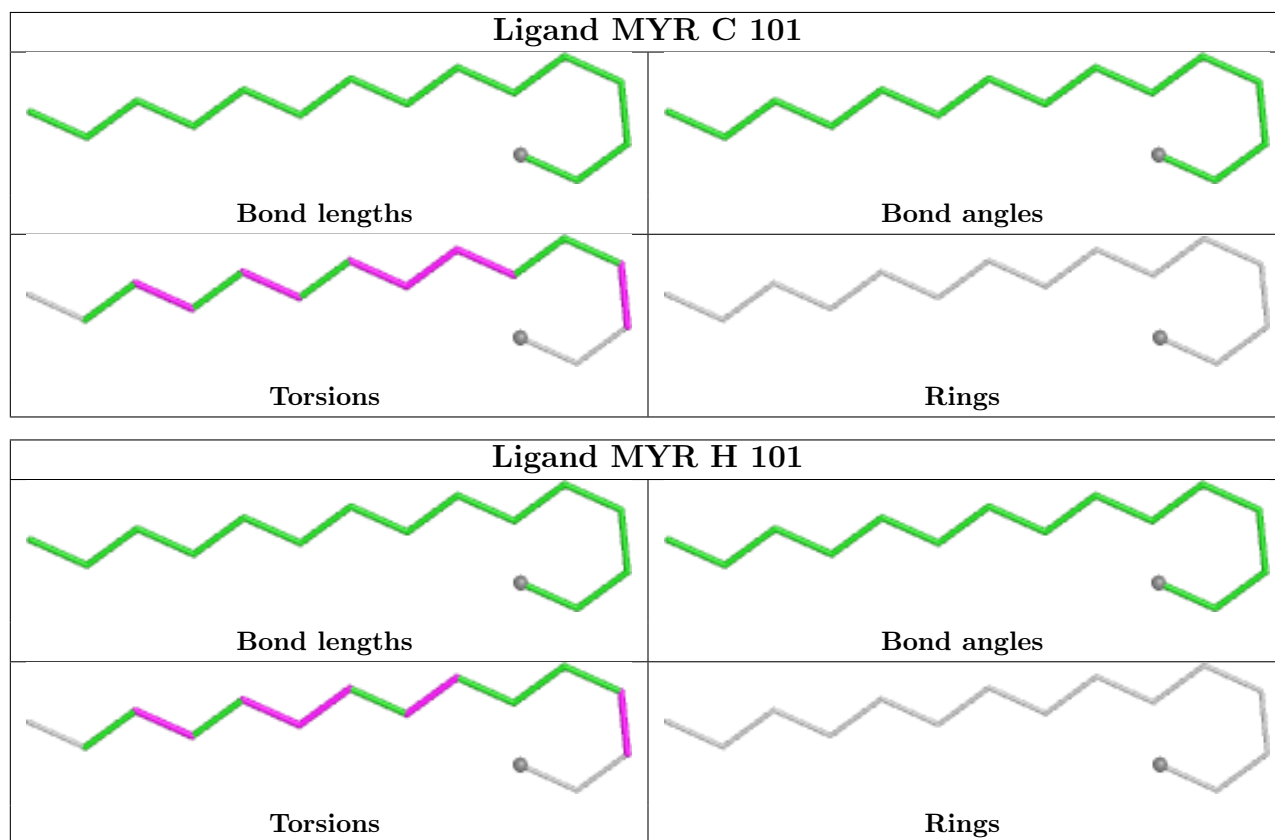
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	J	307	EDO	1	0
7	D	202	EDO	2	0
7	D	203	EDO	1	0
7	A	304	EDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	303	EDO	1	0
9	H	101	MYR	1	0
7	B	106	EDO	1	0
7	J	306	EDO	1	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

There are no chain breaks in this entry.

6 Fit of model and data i

6.1 Protein, DNA and RNA chains i

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	276/276 (100%)	0.11	8 (2%) 53 50	29, 59, 101, 114	1 (0%)
1	F	276/276 (100%)	0.01	7 (2%) 58 55	28, 57, 104, 119	1 (0%)
2	B	100/100 (100%)	-0.30	0 100 100	37, 47, 77, 88	0
2	G	100/100 (100%)	-0.07	1 (1%) 79 78	41, 59, 85, 93	0
3	C	4/4 (100%)	0.06	0 100 100	59, 60, 64, 65	0
3	H	4/4 (100%)	0.80	0 100 100	50, 53, 54, 58	0
4	D	197/197 (100%)	0.59	19 (9%) 13 11	47, 77, 109, 117	0
4	I	197/197 (100%)	-0.11	0 100 100	37, 50, 78, 113	0
5	E	245/245 (100%)	0.40	10 (4%) 41 37	52, 75, 99, 115	0
5	J	245/245 (100%)	-0.17	1 (0%) 88 87	38, 52, 76, 97	0
All	All	1644/1644 (100%)	0.09	46 (2%) 55 51	28, 61, 101, 119	2 (0%)

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	162	SER	3.6
4	D	147	ASP	3.4
1	A	197	HIS	3.1
5	E	220	SER	3.1
5	E	118	ASP	3.1
1	A	200	THR	3.0
4	D	128	THR	3.0
4	D	126	ASN	2.9
5	E	38	ASP	2.8
1	A	199	ALA	2.8
1	F	249	VAL	2.7
4	D	120	LEU	2.7
4	D	139	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
5	E	242	TRP	2.7
1	F	199	ALA	2.6
4	D	195	PHE	2.6
5	E	39	SER	2.6
1	A	257	TYR	2.5
4	D	159	ASP	2.4
1	F	197	HIS	2.4
4	D	164	ASP	2.4
4	D	161	ARG	2.4
4	D	118	TYR	2.3
1	A	201	LEU	2.3
5	E	245	ALA	2.3
4	D	138	SER	2.3
1	A	194	VAL	2.3
1	F	198	GLU	2.2
5	J	62	TYR	2.2
1	F	15	PRO	2.2
4	D	96	LEU	2.2
4	D	140	MET	2.2
4	D	150	VAL	2.2
4	D	95	VAL	2.1
1	F	230	PHE	2.1
5	E	103	TYR	2.1
1	A	1	GLY	2.1
5	E	219	LEU	2.1
1	F	265	GLY	2.1
5	E	102	VAL	2.1
4	D	196	PHE	2.1
4	D	156	CYS	2.0
5	E	129	VAL	2.0
2	G	99	MET	2.0
4	D	197	PRO	2.0
1	A	248	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

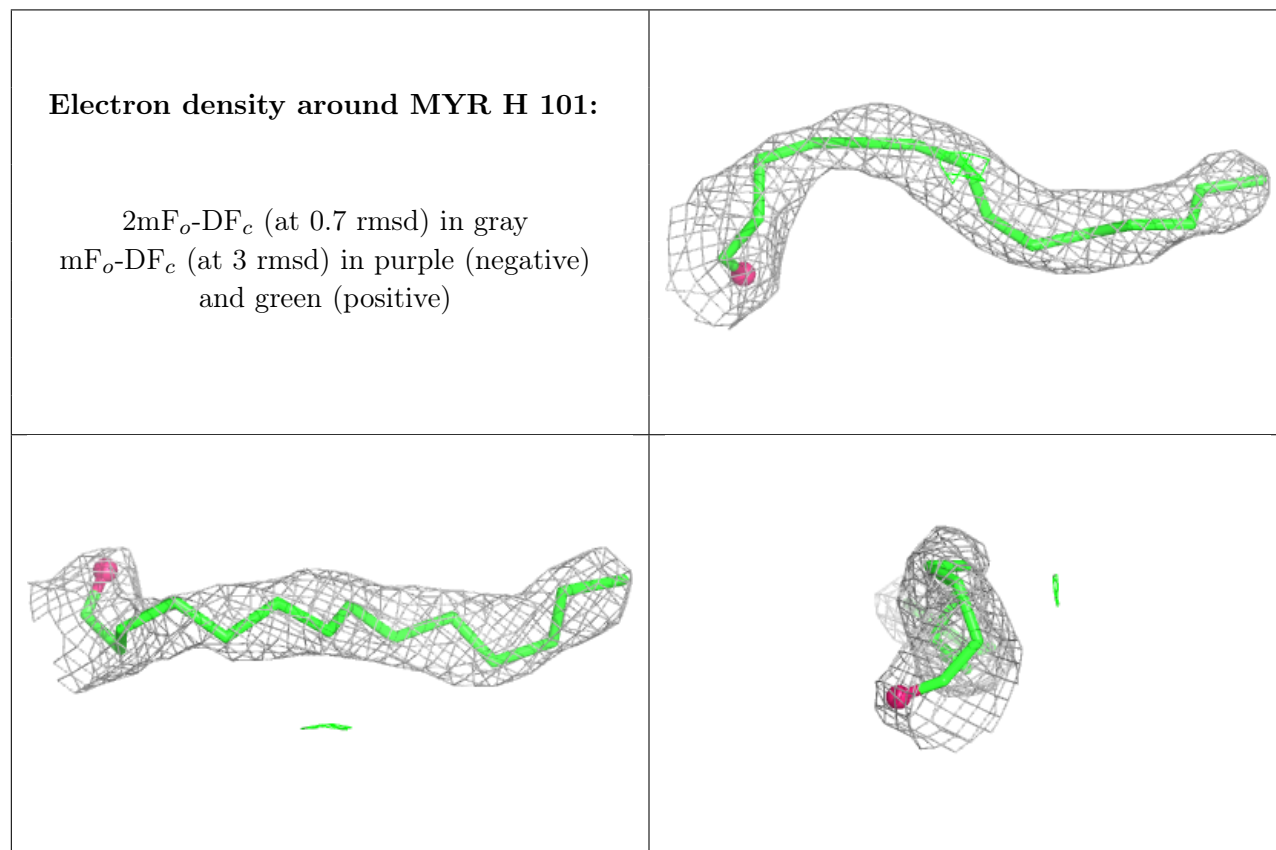
There are no oligosaccharides in this entry.

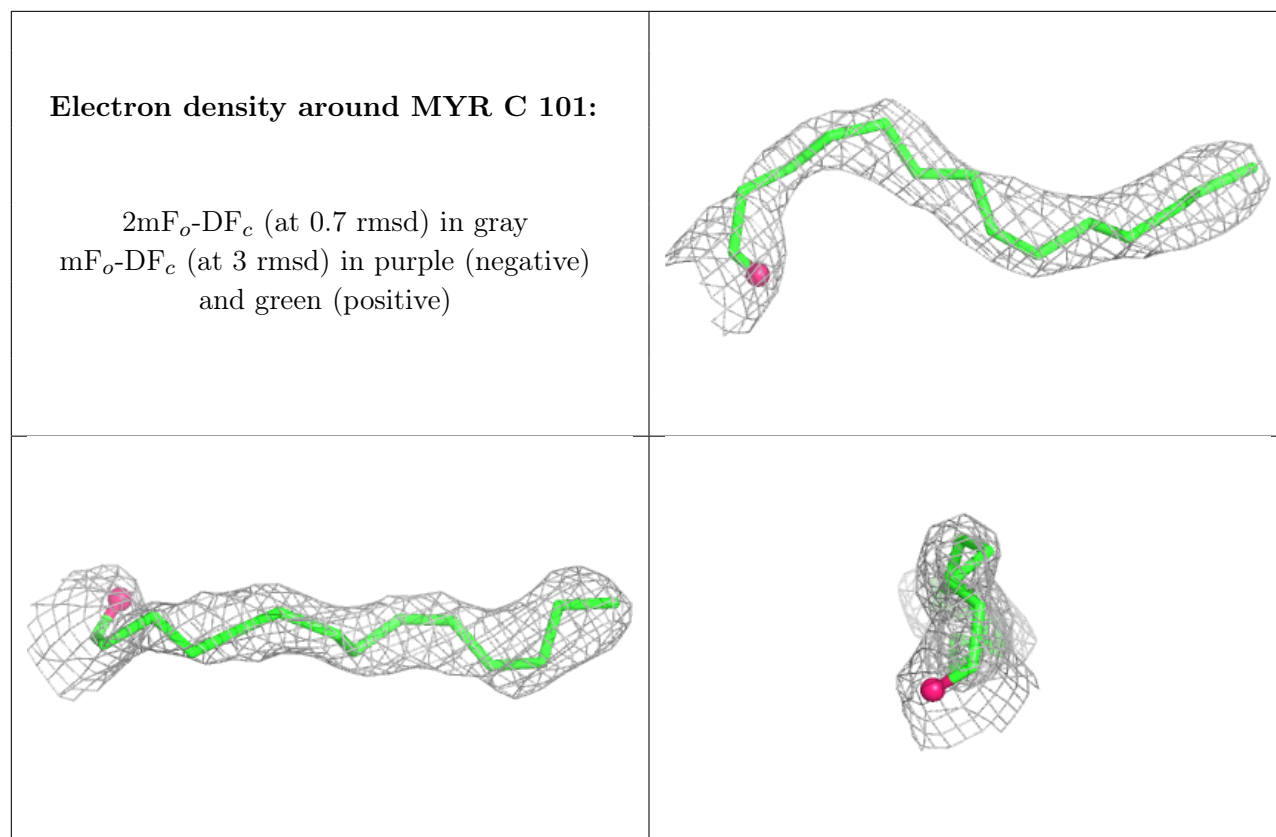
6.4 Ligands i

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	NA	G	102	1/1	0.77	0.20	82,82,82,82	0
7	EDO	A	304	4/4	0.82	0.17	66,67,79,83	0
8	NA	B	103	1/1	0.84	0.09	65,65,65,65	0
7	EDO	B	106	4/4	0.85	0.13	51,57,57,60	0
7	EDO	A	305	4/4	0.85	0.15	55,62,62,71	0
7	EDO	B	105	4/4	0.85	0.12	47,47,49,51	0
7	EDO	D	202	4/4	0.90	0.10	64,69,72,76	0
7	EDO	I	205	4/4	0.90	0.13	52,56,56,58	0
6	IOD	J	303	1/1	0.91	0.13	142,142,142,142	0
7	EDO	A	306	4/4	0.91	0.15	49,57,74,81	0
6	IOD	G	101	1/1	0.92	0.13	122,122,122,122	0
7	EDO	B	104	4/4	0.92	0.10	42,43,47,47	0
7	EDO	D	203	4/4	0.93	0.10	38,45,55,66	0
6	IOD	E	301	1/1	0.93	0.15	161,161,161,161	0
6	IOD	J	305	1/1	0.94	0.15	151,151,151,151	0
6	IOD	F	301	1/1	0.94	0.12	149,149,149,149	0
6	IOD	J	304	1/1	0.94	0.17	133,133,133,133	0
7	EDO	A	303	4/4	0.95	0.08	45,49,50,54	0
8	NA	I	204	1/1	0.95	0.09	61,61,61,61	0
9	MYR	H	101	15/16	0.95	0.10	39,45,59,62	0
6	IOD	E	302	1/1	0.96	0.12	114,114,114,114	0
7	EDO	J	306	4/4	0.96	0.10	51,54,56,59	0
7	EDO	J	307	4/4	0.96	0.08	39,44,45,54	0
6	IOD	B	102	1/1	0.96	0.11	128,128,128,128	0
6	IOD	B	101	1/1	0.96	0.11	111,111,111,111	0
6	IOD	I	201	1/1	0.96	0.18	131,131,131,131	0
9	MYR	C	101	15/16	0.96	0.10	40,53,62,65	0
7	EDO	F	303	4/4	0.96	0.09	43,48,49,50	0
6	IOD	D	201	1/1	0.97	0.21	133,133,133,133	0
6	IOD	A	301	1/1	0.98	0.10	117,117,117,117	0
6	IOD	A	302	1/1	0.98	0.10	102,102,102,102	0
6	IOD	I	202	1/1	0.98	0.19	115,115,115,115	0
6	IOD	J	302	1/1	0.98	0.17	122,122,122,122	0
6	IOD	F	302	1/1	0.98	0.11	97,97,97,97	0
6	IOD	I	203	1/1	0.99	0.09	76,76,76,76	0
6	IOD	J	301	1/1	0.99	0.07	78,78,78,78	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.