



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:36 PM UTC

PDB ID : 8V1L / pdb_00008v1l
Title : Crystal structure of the NTF2L domain of human G3BP1 in complex with small molecule
Authors : Hughes, M.P.; Taylor, J.P.
Deposited on : 2023-11-20
Resolution : 2.68 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

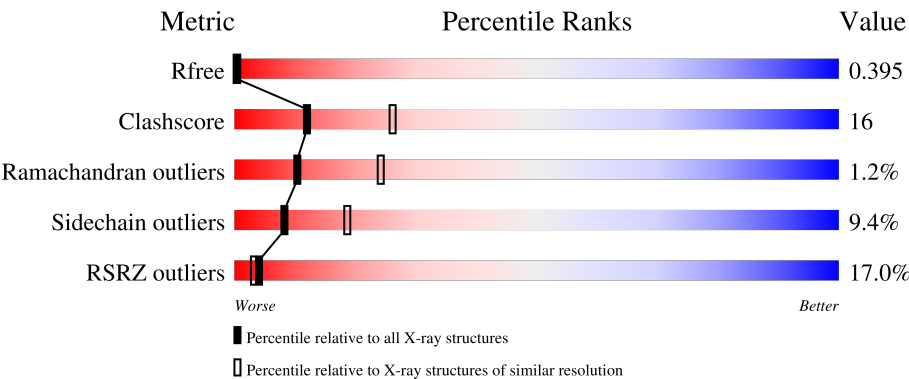
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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	139	17% 50%37%9%
1	B	139	14% 68%26%2%
1	C	139	14% 63%24%9%
1	D	139	9% 59%30%8%
1	E	139	19% 55%37%2%

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Mol	Chain	Length	Quality of chain
1	F	139	22% 65% 24% • 8%

2 Entry composition [i](#)

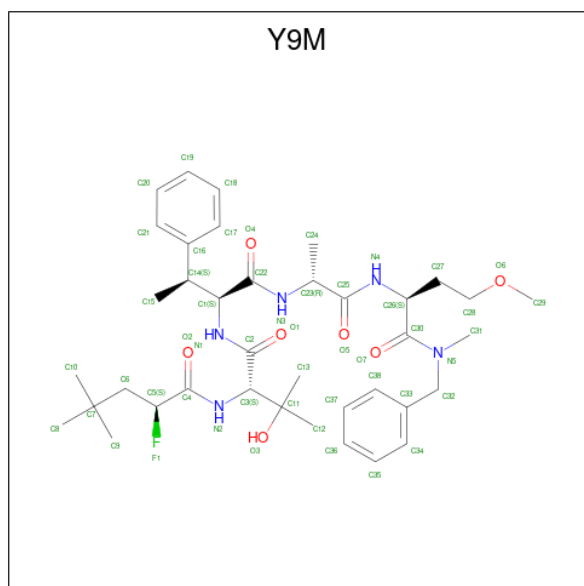
There are 3 unique types of molecules in this entry. The entry contains 6490 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ras GTPase-activating protein-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	126	Total	C	N	O	S	0	1	0
			1004	641	179	178	6			
1	B	133	Total	C	N	O	S	0	1	0
			1044	661	188	190	5			
1	C	127	Total	C	N	O	S	0	2	0
			1010	647	179	179	5			
1	D	128	Total	C	N	O	S	0	0	0
			1013	641	182	185	5			
1	E	134	Total	C	N	O	S	0	2	0
			1023	650	181	187	5			
1	F	128	Total	C	N	O	S	0	2	0
			1013	645	184	178	6			

- Molecule 2 is N-[(2S)-2-fluoro-4,4-dimethylpentanoyl]-3-hydroxy-L-valyl-(betaS)-beta-methyl-L-phenylalanyl-D-alanyl-N-benzyl-N,O-dimethyl-L-homoserinamide (CCD ID: Y9M) (formula: C₃₈H₅₆FN₅O₇) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			51	38	1	5	7		
2	B	1	Total	C	F	N	O	0	0
			48	36	1	5	6		
2	C	1	Total	C	F	N	O	0	0
			45	32	1	5	7		
2	D	1	Total	C	F	N	O	0	0
			51	38	1	5	7		
2	E	1	Total	C	F	N	O	0	0
			45	32	1	5	7		
2	F	1	Total	C	F	N	O	0	0
			51	38	1	5	7		

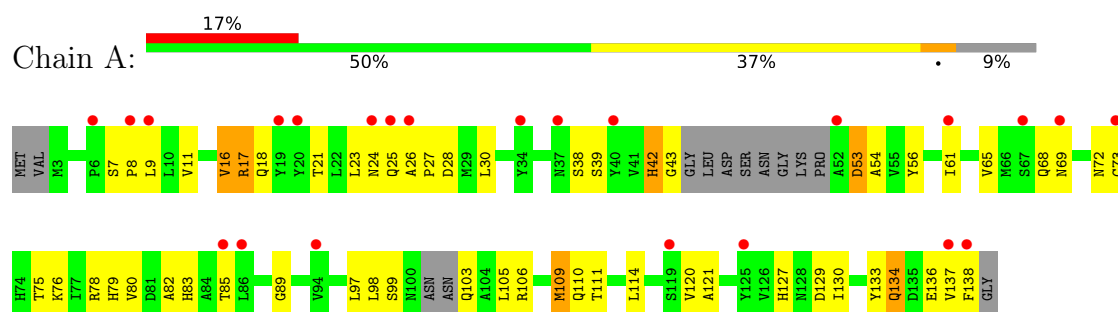
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total	O	0	0
			6	6		
3	B	10	Total	O	0	0
			10	10		
3	C	11	Total	O	0	0
			11	11		
3	D	16	Total	O	0	0
			16	16		
3	E	21	Total	O	0	0
			21	21		
3	F	28	Total	O	0	0
			28	28		

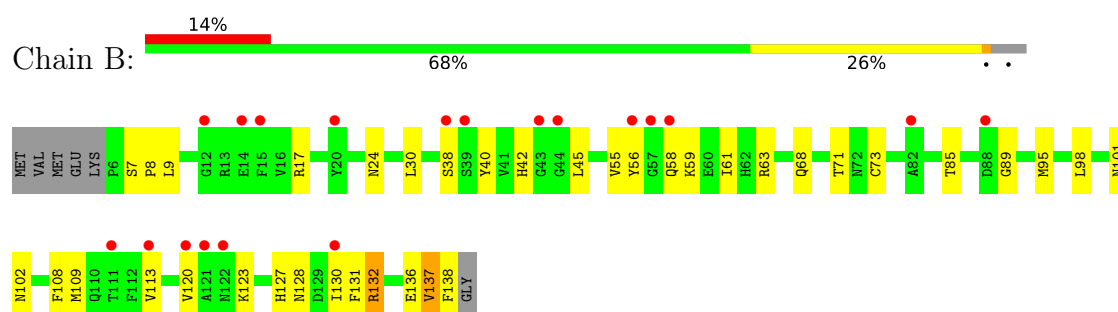
3 Residue-property plots [i](#)

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.

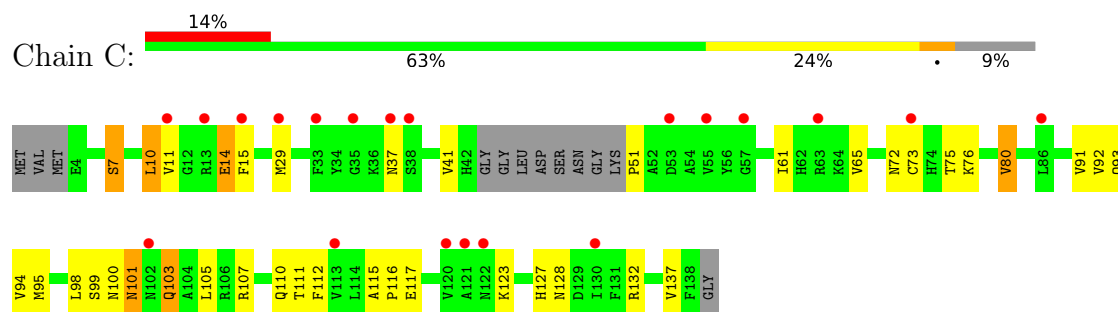
- Molecule 1: Ras GTPase-activating protein-binding protein 1



- Molecule 1: Ras GTPase-activating protein-binding protein 1

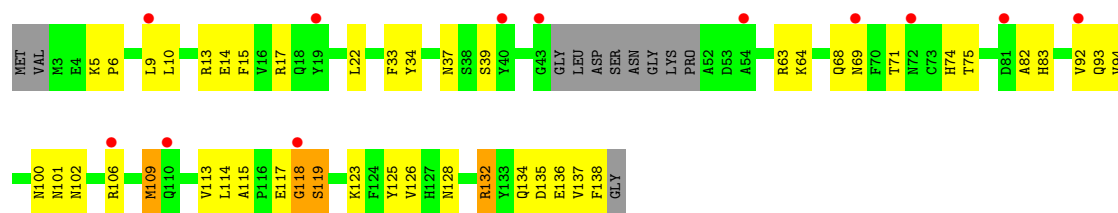


- Molecule 1: Ras GTPase-activating protein-binding protein 1

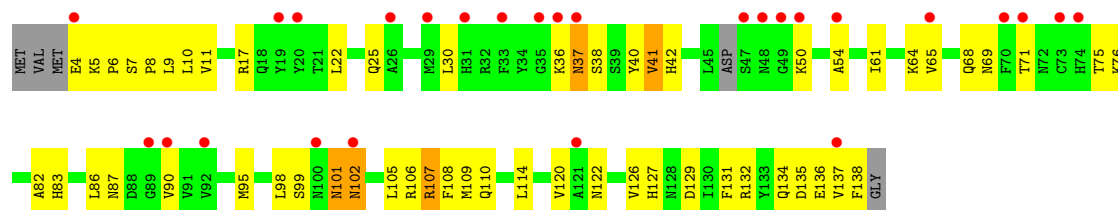


- Molecule 1: Ras GTPase-activating protein-binding protein 1

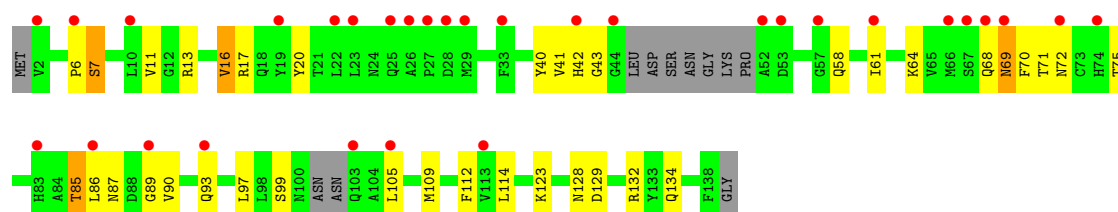




● Molecule 1: Ras GTPase-activating protein-binding protein 1



● Molecule 1: Ras GTPase-activating protein-binding protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.17Å 84.69Å 102.15Å 90.00° 91.26° 90.00°	Depositor
Resolution (Å)	46.87 – 2.68 46.87 – 2.68	Depositor EDS
% Data completeness (in resolution range)	96.5 (46.87-2.68) 96.5 (46.87-2.68)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.276 , 0.395 0.278 , 0.395	Depositor DCC
R_{free} test set	1165 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	39.6	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	6490	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 31.35 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1232e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: Y9M

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	0.48	0/1030	0.97	0/1394
1	B	0.49	0/1073	0.94	0/1455
1	C	0.50	0/1041	0.94	0/1413
1	D	0.50	0/1037	0.94	0/1407
1	E	0.51	0/1054	0.97	0/1432
1	F	0.49	0/1039	0.92	0/1405
All	All	0.49	0/6274	0.95	0/8506

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
1	A	0	2
1	B	0	1
1	D	0	1
1	E	0	1
1	F	0	1
All	All	0	6

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	17	ARG	Sidechain
1	A	78	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	B	17	ARG	Sidechain
1	D	63	ARG	Sidechain
1	E	50	LYS	Peptide
1	F	13	ARG	Sidechain

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	1004	0	949	42	0
1	B	1044	0	978	36	0
1	C	1010	0	956	33	0
1	D	1013	0	940	38	0
1	E	1023	0	928	33	1
1	F	1013	0	954	23	1
2	A	51	0	0	1	0
2	B	48	0	0	0	0
2	C	45	0	0	3	0
2	D	51	0	0	1	0
2	E	45	0	0	1	0
2	F	51	0	0	3	0
3	A	6	0	0	2	0
3	B	10	0	0	0	0
3	C	11	0	0	1	0
3	D	16	0	0	2	0
3	E	21	0	0	3	0
3	F	28	0	0	3	0
All	All	6490	0	5705	190	1

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 16.

All (190) 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:F:17:ARG:NH2	3:F:301:HOH:O	1.93	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:MET:HE1	1:D:134:GLN:HA	1.58	0.86
1:C:103:GLN:HE21	1:C:103:GLN:HA	1.40	0.85
1:E:6:PRO:CD	1:E:122:ASN:HD21	1.97	0.78
1:A:99:SER:HB2	1:A:103:GLN:O	1.84	0.77
1:A:42:HIS:CG	1:A:68:GLN:HE22	2.05	0.74
1:D:14:GLU:OE1	1:D:17:ARG:NH2	2.20	0.73
1:D:137:VAL:HG12	1:D:137:VAL:O	1.89	0.71
1:D:137:VAL:O	1:D:137:VAL:CG1	2.41	0.69
1:C:103:GLN:HE21	1:C:103:GLN:CA	2.06	0.68
1:F:93:GLN:O	3:F:302:HOH:O	2.11	0.68
1:F:7:SER:O	1:F:11:VAL:HG23	1.94	0.68
1:A:61:ILE:O	1:A:65:VAL:HG23	1.94	0.66
1:C:101:ASN:OD1	1:E:25:GLN:NE2	2.29	0.66
1:A:109:MET:HE2	1:C:93:GLN:HG3	1.77	0.66
1:D:9:LEU:HB2	3:D:312:HOH:O	1.98	0.64
1:A:99:SER:CB	1:A:103:GLN:O	2.46	0.62
1:B:85:THR:HG22	1:B:89:GLY:C	2.25	0.61
1:E:22:LEU:HD23	1:E:30:LEU:HD12	1.82	0.61
1:C:103:GLN:HA	1:C:103:GLN:NE2	2.14	0.61
1:E:7:SER:O	1:E:11:VAL:HG23	2.00	0.61
1:A:130:ILE:HG22	1:C:93:GLN:NE2	2.15	0.61
1:B:40:TYR:HB3	1:B:61:ILE:HG12	1.84	0.60
1:D:64:LYS:C	1:D:68:GLN:HE21	2.09	0.60
1:E:10:LEU:CB	3:E:310:HOH:O	2.50	0.60
1:A:42:HIS:O	1:A:43:GLY:C	2.44	0.59
1:D:115:ALA:HB3	1:D:125:TYR:CE1	2.38	0.59
1:A:97:LEU:HB3	1:A:105:LEU:HB3	1.85	0.59
1:F:61:ILE:HG23	3:F:308:HOH:O	2.03	0.59
1:B:71:THR:OG1	1:B:101:ASN:OD1	2.21	0.58
1:D:132:ARG:NH1	1:D:136:GLU:OE1	2.37	0.57
1:E:110:GLN:NE2	1:E:129:ASP:OD1	2.37	0.57
1:E:37:ASN:ND2	1:E:127:HIS:CE1	2.73	0.57
1:C:123:LYS:HG3	2:C:201:Y9M:O3	2.05	0.56
1:A:26:ALA:HA	3:A:302:HOH:O	2.06	0.56
1:E:95:MET:HE3	1:E:109:MET:HE2	1.85	0.56
1:E:101:ASN:O	1:E:102:ASN:CB	2.55	0.55
1:D:22:LEU:HD23	1:D:33:PHE:CZ	2.42	0.55
1:D:132:ARG:HH11	1:D:132:ARG:HA	1.70	0.55
1:A:134:GLN:HA	1:C:95:MET:CE	2.38	0.54
1:B:95:MET:HE3	1:B:109:MET:CE	2.37	0.54
1:D:134:GLN:HG2	1:D:138:PHE:CD2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:GLN:HA	1:C:95:MET:HE1	1.89	0.54
1:D:123:LYS:NZ	2:D:201:Y9M:O5	2.25	0.54
1:E:38:SER:HA	1:E:127:HIS:O	2.08	0.54
1:A:79:HIS:CD2	1:A:80:VAL:N	2.76	0.53
1:C:98[A]:LEU:HD23	1:C:99:SER:N	2.24	0.53
1:C:98[B]:LEU:HD23	1:C:99:SER:N	2.24	0.53
1:E:83:HIS:O	1:E:90:VAL:HG13	2.08	0.53
1:E:114:LEU:CD2	1:E:126:VAL:HG22	2.39	0.53
1:A:138:PHE:CE2	1:C:107:ARG:HG2	2.44	0.53
1:B:95:MET:HE2	1:D:138:PHE:HE2	1.74	0.53
1:D:82:ALA:C	1:D:83:HIS:ND1	2.68	0.52
1:F:85:THR:HG22	1:F:89:GLY:C	2.34	0.52
1:E:17:ARG:HG2	1:E:17:ARG:HH11	1.74	0.51
1:B:101:ASN:O	1:B:102:ASN:C	2.53	0.51
1:E:98[A]:LEU:HD23	1:E:99:SER:N	2.26	0.51
1:E:98[B]:LEU:HD23	1:E:99:SER:N	2.26	0.51
1:B:42:HIS:O	1:B:131:PHE:O	2.28	0.51
1:D:74:HIS:NE2	1:D:102:ASN:ND2	2.59	0.51
1:F:68:GLN:O	1:F:70:PHE:N	2.43	0.51
1:A:111:THR:HG23	1:C:111:THR:HG21	1.93	0.51
1:B:38:SER:HA	1:B:127:HIS:O	2.11	0.51
1:B:132:ARG:HH11	1:B:132:ARG:HG3	1.75	0.51
1:E:61:ILE:O	1:E:65:VAL:HG23	2.11	0.51
1:A:39:SER:HB2	1:A:54:ALA:HB1	1.92	0.51
1:B:55:VAL:C	1:B:56:TYR:CD1	2.90	0.50
1:E:41:VAL:HG22	1:E:54:ALA:HB2	1.93	0.50
1:B:85:THR:OG1	1:D:113:VAL:HG11	2.11	0.50
1:A:103:GLN:OE1	1:F:132:ARG:NH2	2.43	0.50
1:C:123:LYS:NZ	2:C:201:Y9M:O5	2.32	0.50
2:E:201:Y9M:O6	2:E:201:Y9M:N4	2.44	0.50
1:F:85:THR:HG22	1:F:89:GLY:CA	2.42	0.50
1:B:120:VAL:O	1:B:123:LYS:HB3	2.12	0.50
1:C:7:SER:O	1:C:11:VAL:HG23	2.12	0.49
1:F:123:LYS:NZ	2:F:201:Y9M:O5	2.44	0.49
1:A:9:LEU:HA	1:A:82:ALA:HB3	1.94	0.49
1:B:95:MET:HE3	1:B:109:MET:HE2	1.93	0.49
1:E:6:PRO:HD3	1:E:122:ASN:HD21	1.75	0.49
1:D:5:LYS:O	3:D:301:HOH:O	2.18	0.49
1:C:73:CYS:SG	1:C:98[B]:LEU:HD21	2.53	0.48
1:A:85:THR:HA	1:C:127:HIS:CD2	2.49	0.48
1:D:100:ASN:O	1:D:101:ASN:C	2.57	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:GLN:NE2	1:A:129:ASP:OD1	2.41	0.48
1:B:130:ILE:HG22	1:D:93:GLN:NE2	2.29	0.47
1:C:73:CYS:SG	1:C:98[A]:LEU:HD21	2.53	0.47
1:B:73:CYS:SG	1:B:98[B]:LEU:HD21	2.54	0.47
1:C:103:GLN:CA	1:C:103:GLN:NE2	2.77	0.47
1:E:42:HIS:HA	1:E:131:PHE:HB3	1.96	0.47
1:B:73:CYS:SG	1:B:98[A]:LEU:HD21	2.54	0.47
1:E:6:PRO:HD2	1:E:122:ASN:HD21	1.79	0.47
1:F:112:PHE:HA	1:F:128:ASN:O	2.15	0.47
1:D:132:ARG:NH2	1:D:136:GLU:OE1	2.48	0.47
1:E:137:VAL:HG12	1:E:138:PHE:N	2.30	0.47
1:A:82:ALA:C	1:A:83:HIS:HD1	2.23	0.47
1:B:24:ASN:ND2	1:B:73:CYS:O	2.39	0.46
1:B:113:VAL:HB	1:B:128:ASN:HB2	1.97	0.46
1:C:100:ASN:O	1:C:101:ASN:C	2.58	0.46
1:F:58:GLN:HE22	2:F:201:Y9M:C29	2.28	0.46
1:F:71:THR:O	1:F:72:ASN:C	2.58	0.46
1:A:106:ARG:HB2	1:A:133:TYR:CE1	2.51	0.46
2:F:201:Y9M:O6	2:F:201:Y9M:N4	2.49	0.46
1:B:45:LEU:HD21	1:B:136:GLU:HB3	1.98	0.46
1:C:61:ILE:O	1:C:65:VAL:HG23	2.15	0.46
1:A:134:GLN:O	1:A:137:VAL:O	2.34	0.46
1:B:109:MET:SD	1:D:109:MET:HB3	2.56	0.46
1:A:127:HIS:HD2	1:A:127:HIS:O	2.00	0.45
1:D:94:VAL:O	1:D:109:MET:HA	2.16	0.45
1:E:8:PRO:HA	1:E:11:VAL:HB	1.98	0.45
1:B:137:VAL:O	1:B:138:PHE:C	2.59	0.45
1:A:75:THR:HG22	1:A:76:LYS:H	1.82	0.45
1:C:115:ALA:O	1:C:116:PRO:C	2.59	0.45
1:B:42:HIS:CB	1:B:68:GLN:HE22	2.29	0.45
1:D:113:VAL:HG23	1:D:128:ASN:HB2	1.98	0.45
1:B:108:PHE:CG	1:B:131:PHE:CE2	3.05	0.45
1:A:127:HIS:O	1:A:127:HIS:CD2	2.70	0.45
1:D:14:GLU:O	1:D:15:PHE:C	2.58	0.45
1:F:68:GLN:O	1:F:69:ASN:C	2.59	0.45
1:F:42[B]:HIS:CG	1:F:68:GLN:OE1	2.70	0.44
1:D:118:GLY:O	1:D:119:SER:HB3	2.16	0.44
1:E:122:ASN:HD22	1:E:122:ASN:HA	1.53	0.44
1:F:16:VAL:O	1:F:20:TYR:HD2	2.00	0.44
1:F:97:LEU:HB3	1:F:105:LEU:HB3	1.99	0.44
1:A:53:ASP:OD1	1:A:53:ASP:N	2.43	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:114:LEU:CD2	1:D:126:VAL:HG22	2.47	0.44
1:B:8:PRO:O	1:B:9:LEU:C	2.61	0.44
1:D:22:LEU:HD23	1:D:33:PHE:CE2	2.53	0.44
1:A:85:THR:HG22	1:A:89:GLY:C	2.43	0.44
1:A:16:VAL:O	1:A:17:ARG:C	2.61	0.44
1:E:40:TYR:CB	1:E:61:ILE:HG12	2.48	0.44
1:F:40:TYR:CE1	1:F:64:LYS:HG3	2.53	0.44
1:C:14:GLU:HG2	1:D:10:LEU:HD23	2.00	0.43
1:A:98[A]:LEU:HD23	1:A:133:TYR:OH	2.18	0.43
1:E:68:GLN:O	1:E:69:ASN:C	2.61	0.43
1:C:7:SER:CB	1:C:10:LEU:HD22	2.48	0.43
1:B:7:SER:HA	1:B:8:PRO:HD3	1.89	0.43
1:B:42:HIS:HB3	1:B:68:GLN:HE22	1.84	0.43
1:C:72:ASN:ND2	3:C:304:HOH:O	2.51	0.43
1:D:109:MET:O	1:D:109:MET:HG3	2.18	0.43
1:E:9:LEU:HB2	3:E:314:HOH:O	2.16	0.43
1:A:24:ASN:ND2	1:A:73:CYS:O	2.48	0.43
1:B:58:GLN:O	1:B:59:LYS:C	2.60	0.43
1:A:27:PRO:HD2	3:A:302:HOH:O	2.19	0.43
1:A:120:VAL:HG12	1:A:121:ALA:O	2.19	0.43
1:D:109:MET:HE3	1:D:109:MET:HB2	1.85	0.43
1:B:56:TYR:CD1	1:B:56:TYR:N	2.87	0.43
1:E:134:GLN:O	1:E:138:PHE:HB2	2.18	0.43
1:F:6:PRO:O	1:F:7:SER:C	2.61	0.43
1:A:25:GLN:O	1:A:26:ALA:C	2.61	0.42
1:C:80:VAL:HG22	1:C:94:VAL:HG22	2.01	0.42
1:B:95:MET:CE	1:D:134:GLN:HA	2.40	0.42
1:F:99:SER:CB	1:F:105:LEU:HD23	2.50	0.42
1:F:90:VAL:HB	1:F:114:LEU:HB2	2.00	0.42
1:B:95:MET:HE3	1:B:109:MET:HE3	2.02	0.42
1:D:117:GLU:O	1:D:118:GLY:C	2.62	0.42
1:E:108:PHE:C	1:E:108:PHE:CD1	2.96	0.42
1:B:132:ARG:HH12	1:B:136:GLU:HB2	1.84	0.42
1:D:68:GLN:O	1:D:69:ASN:C	2.62	0.42
1:A:137:VAL:HG12	1:A:138:PHE:N	2.35	0.42
1:E:40:TYR:HB2	1:E:61:ILE:HG12	2.02	0.42
1:B:95:MET:HB2	1:D:137:VAL:HG11	2.02	0.42
1:A:7:SER:O	1:A:11:VAL:HG23	2.19	0.42
1:C:101:ASN:O	1:C:103:GLN:HG2	2.20	0.42
1:C:92:VAL:HG12	1:C:94:VAL:HG23	2.01	0.42
1:E:86:LEU:O	1:E:87:ASN:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:85:THR:HG22	1:F:89:GLY:HA3	2.01	0.42
1:F:86:LEU:O	1:F:87:ASN:C	2.63	0.42
1:D:92:VAL:HG12	1:D:94:VAL:HG23	2.02	0.41
1:A:98[B]:LEU:HD23	1:A:133:TYR:OH	2.20	0.41
1:B:40:TYR:CD1	1:B:40:TYR:C	2.97	0.41
1:A:42:HIS:CB	1:A:68:GLN:HE22	2.32	0.41
1:E:75:THR:OG1	3:E:301:HOH:O	2.04	0.41
1:A:114:LEU:HD13	2:A:201:Y9M:C20	2.50	0.41
1:E:9:LEU:HA	1:E:82:ALA:HB3	2.02	0.41
1:C:15:PHE:HA	2:C:201:Y9M:C18	2.51	0.41
1:C:99:SER:CB	1:C:105:LEU:HD23	2.51	0.41
1:D:33:PHE:HB2	1:D:34:TYR:CE1	2.55	0.41
1:A:21:THR:C	1:A:23:LEU:H	2.29	0.41
1:C:112:PHE:HA	1:C:128:ASN:O	2.20	0.41
1:C:110:GLN:CD	1:C:112:PHE:CE2	2.99	0.40
1:A:38:SER:O	1:A:56:TYR:HA	2.22	0.40
1:C:98[B]:LEU:HD23	1:C:99:SER:O	2.21	0.40
1:F:134:GLN:HE21	1:F:134:GLN:HB3	1.71	0.40
1:D:17:ARG:HG2	1:D:17:ARG:HH11	1.86	0.40
1:E:64:LYS:O	1:E:68:GLN:HG3	2.21	0.40
1:A:68:GLN:O	1:A:69:ASN:C	2.64	0.40
1:A:134:GLN:C	1:A:136:GLU:H	2.30	0.40
1:B:73:CYS:SG	1:B:98[A]:LEU:HD11	2.61	0.40
1:B:73:CYS:SG	1:B:98[B]:LEU:HD11	2.61	0.40
1:D:114:LEU:HD21	1:D:126:VAL:HG22	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:107:ARG:NH1	1:F:134:GLN:OE1[2_446]	2.06	0.14

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	121/139 (87%)	106 (88%)	14 (12%)	1 (1%)	16	34
1	B	132/139 (95%)	115 (87%)	17 (13%)	0	100	100
1	C	125/139 (90%)	113 (90%)	10 (8%)	2 (2%)	7	18
1	D	124/139 (89%)	109 (88%)	12 (10%)	3 (2%)	4	10
1	E	132/139 (95%)	117 (89%)	14 (11%)	1 (1%)	16	34
1	F	124/139 (89%)	109 (88%)	13 (10%)	2 (2%)	7	18
All	All	758/834 (91%)	669 (88%)	80 (11%)	9 (1%)	10	24

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	119	SER
1	E	102	ASN
1	F	69	ASN
1	D	118	GLY
1	C	101	ASN
1	A	8	PRO
1	F	43	GLY
1	C	80	VAL
1	D	6	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	104/121 (86%)	95 (91%)	9 (9%)	9	21
1	B	108/121 (89%)	104 (96%)	4 (4%)	30	56
1	C	105/121 (87%)	91 (87%)	14 (13%)	4	8
1	D	105/121 (87%)	96 (91%)	9 (9%)	10	22
1	E	101/121 (84%)	86 (85%)	15 (15%)	3	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	103/121 (85%)	96 (93%)	7 (7%)	14	32
All	All	626/726 (86%)	568 (91%)	58 (9%)	8	19

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	VAL
1	A	18	GLN
1	A	28	ASP
1	A	30	LEU
1	A	42	HIS
1	A	53	ASP
1	A	72	ASN
1	A	109	MET
1	A	134	GLN
1	B	30	LEU
1	B	63	ARG
1	B	132	ARG
1	B	137	VAL
1	C	7	SER
1	C	10	LEU
1	C	14	GLU
1	C	29	MET
1	C	37	ASN
1	C	41	VAL
1	C	51	PRO
1	C	75	THR
1	C	76	LYS
1	C	91	VAL
1	C	103	GLN
1	C	117	GLU
1	C	132	ARG
1	C	137	VAL
1	D	13	ARG
1	D	37	ASN
1	D	39	SER
1	D	71	THR
1	D	75	THR
1	D	106	ARG
1	D	109	MET
1	D	132	ARG
1	D	135	ASP

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Mol	Chain	Res	Type
1	E	4	GLU
1	E	5	LYS
1	E	36	LYS
1	E	37	ASN
1	E	41	VAL
1	E	71	THR
1	E	76	LYS
1	E	101	ASN
1	E	105	LEU
1	E	106	ARG
1	E	107	ARG
1	E	120	VAL
1	E	132	ARG
1	E	135	ASP
1	E	136	GLU
1	F	7	SER
1	F	16	VAL
1	F	41	VAL
1	F	75	THR
1	F	85	THR
1	F	109	MET
1	F	129	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	GLN
1	A	68	GLN
1	A	100	ASN
1	A	127	HIS
1	B	58	GLN
1	B	62	HIS
1	B	68	GLN
1	B	69	ASN
1	B	79	HIS
1	C	37	ASN
1	C	58	GLN
1	C	62	HIS
1	C	72	ASN
1	C	103	GLN
1	C	127	HIS
1	C	128	ASN

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Mol	Chain	Res	Type
1	D	37	ASN
1	D	58	GLN
1	D	68	GLN
1	D	72	ASN
1	D	102	ASN
1	D	127	HIS
1	E	37	ASN
1	E	68	GLN
1	E	122	ASN
1	E	127	HIS
1	E	128	ASN
1	F	58	GLN
1	F	74	HIS
1	F	93	GLN
1	F	122	ASN
1	F	134	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 [i](#)

6 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	Y9M	F	201	-	49,52,52	0.28	0	60,73,73	0.49	0
2	Y9M	B	201	-	46,49,52	0.28	0	58,70,73	0.93	2 (3%)
2	Y9M	C	201	-	42,45,52	0.38	0	51,64,73	0.83	2 (3%)
2	Y9M	A	201	-	49,52,52	0.30	0	60,73,73	0.84	1 (1%)
2	Y9M	D	201	-	49,52,52	0.28	0	60,73,73	0.90	4 (6%)
2	Y9M	E	201	-	42,45,52	0.27	0	51,64,73	0.56	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
2	Y9M	F	201	-	-	13/67/67/67	0/2/2/2
2	Y9M	B	201	-	-	13/63/63/67	0/2/2/2
2	Y9M	C	201	-	-	14/63/63/67	0/1/1/2
2	Y9M	A	201	-	-	6/67/67/67	0/2/2/2
2	Y9M	D	201	-	-	6/67/67/67	0/2/2/2
2	Y9M	E	201	-	-	8/63/63/67	0/1/1/2

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	Y9M	C7-C6-C5	-5.00	110.12	116.46
2	A	201	Y9M	C7-C6-C5	-4.97	110.15	116.46
2	D	201	Y9M	C7-C6-C5	-3.60	111.89	116.46
2	D	201	Y9M	C12-C11-C3	-3.05	106.19	111.02
2	C	201	Y9M	C13-C11-C3	3.00	115.77	111.02
2	C	201	Y9M	C26-C30-N5	2.35	121.19	118.53
2	B	201	Y9M	C12-C11-C3	-2.08	107.73	111.02
2	D	201	Y9M	C27-C26-C30	2.03	114.79	109.75
2	D	201	Y9M	C1-N1-C2	2.01	126.88	121.90

There are no chirality outliers.

All (60) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	201	Y9M	C26-C27-C28-O6

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Mol	Chain	Res	Type	Atoms
2	B	201	Y9M	C13-C11-C3-C2
2	B	201	Y9M	C13-C11-C3-N2
2	B	201	Y9M	O3-C11-C3-N2
2	C	201	Y9M	N2-C4-C5-F1
2	C	201	Y9M	O2-C4-C5-F1
2	C	201	Y9M	F1-C5-C6-C7
2	C	201	Y9M	N4-C26-C27-C28
2	C	201	Y9M	C30-C26-C27-C28
2	C	201	Y9M	C13-C11-C3-C2
2	D	201	Y9M	O2-C4-C5-F1
2	D	201	Y9M	N4-C26-C27-C28
2	D	201	Y9M	C30-C26-C27-C28
2	E	201	Y9M	C5-C4-N2-C3
2	E	201	Y9M	F1-C5-C6-C7
2	F	201	Y9M	N2-C4-C5-F1
2	F	201	Y9M	O2-C4-C5-F1
2	F	201	Y9M	C5-C6-C7-C8
2	F	201	Y9M	N1-C1-C14-C16
2	E	201	Y9M	O2-C4-N2-C3
2	E	201	Y9M	C23-C25-N4-C26
2	E	201	Y9M	O5-C25-N4-C26
2	B	201	Y9M	C5-C4-N2-C3
2	A	201	Y9M	C27-C28-O6-C29
2	E	201	Y9M	N4-C26-C27-C28
2	F	201	Y9M	N4-C26-C27-C28
2	C	201	Y9M	C27-C28-O6-C29
2	D	201	Y9M	C27-C28-O6-C29
2	F	201	Y9M	C5-C6-C7-C10
2	F	201	Y9M	C5-C6-C7-C9
2	A	201	Y9M	C5-C4-N2-C3
2	D	201	Y9M	C5-C4-N2-C3
2	B	201	Y9M	O2-C4-N2-C3
2	C	201	Y9M	O2-C4-N2-C3
2	B	201	Y9M	O7-C30-N5-C32
2	B	201	Y9M	C12-C11-C3-C2
2	F	201	Y9M	C30-C26-C27-C28
2	F	201	Y9M	C24-C23-C25-O5
2	B	201	Y9M	O3-C11-C3-C2
2	E	201	Y9M	C11-C3-N2-C4
2	C	201	Y9M	C5-C4-N2-C3
2	B	201	Y9M	N3-C23-C25-O5
2	F	201	Y9M	N3-C23-C25-O5

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Mol	Chain	Res	Type	Atoms
2	C	201	Y9M	O1-C2-C3-N2
2	C	201	Y9M	N1-C2-C3-N2
2	B	201	Y9M	N3-C23-C25-N4
2	F	201	Y9M	N3-C23-C25-N4
2	B	201	Y9M	C24-C23-C25-O5
2	F	201	Y9M	C24-C23-C25-N4
2	C	201	Y9M	C27-C26-C30-N5
2	A	201	Y9M	C27-C26-C30-O7
2	A	201	Y9M	N5-C32-C33-C38
2	F	201	Y9M	C27-C26-C30-O7
2	E	201	Y9M	C30-C26-C27-C28
2	B	201	Y9M	C12-C11-C3-N2
2	C	201	Y9M	C13-C11-C3-N2
2	B	201	Y9M	C24-C23-C25-N4
2	C	201	Y9M	C5-C6-C7-C10
2	D	201	Y9M	C27-C26-C30-O7
2	A	201	Y9M	N5-C32-C33-C34

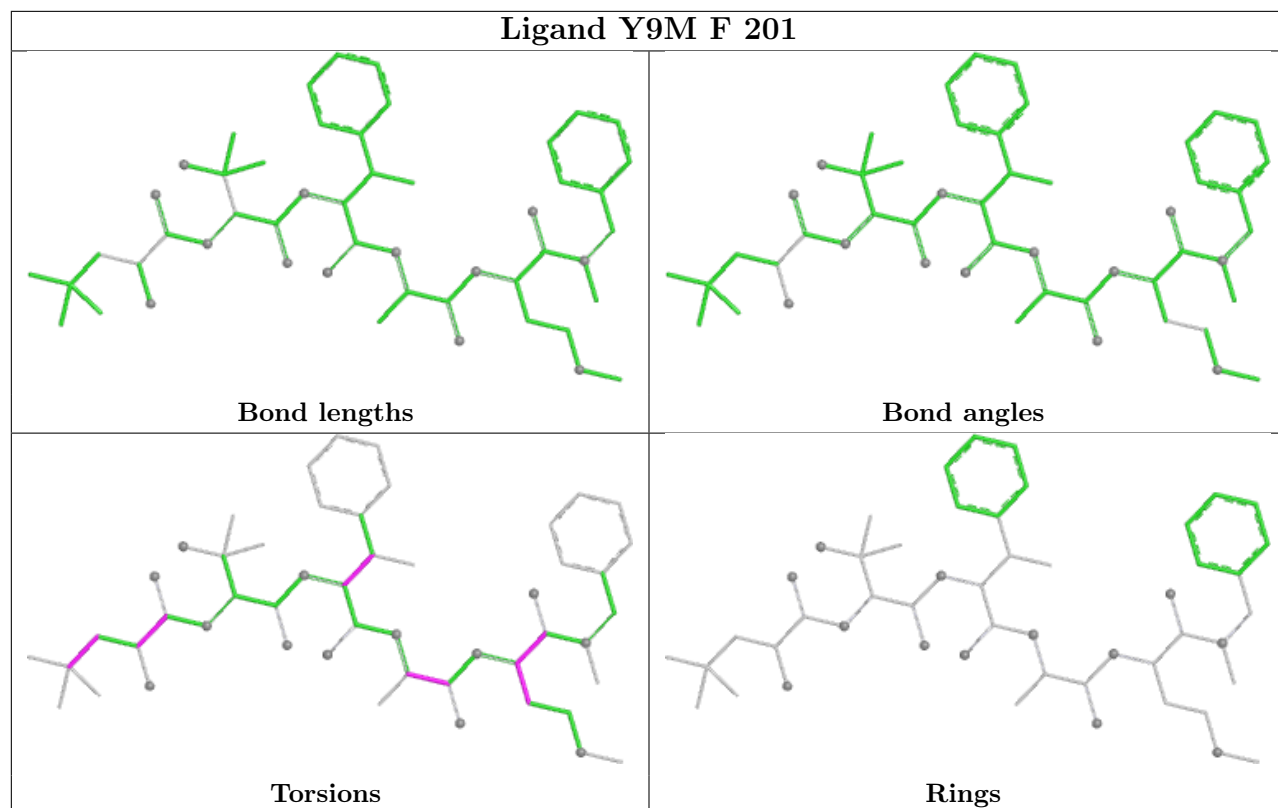
There are no ring outliers.

5 monomers are involved in 9 short contacts:

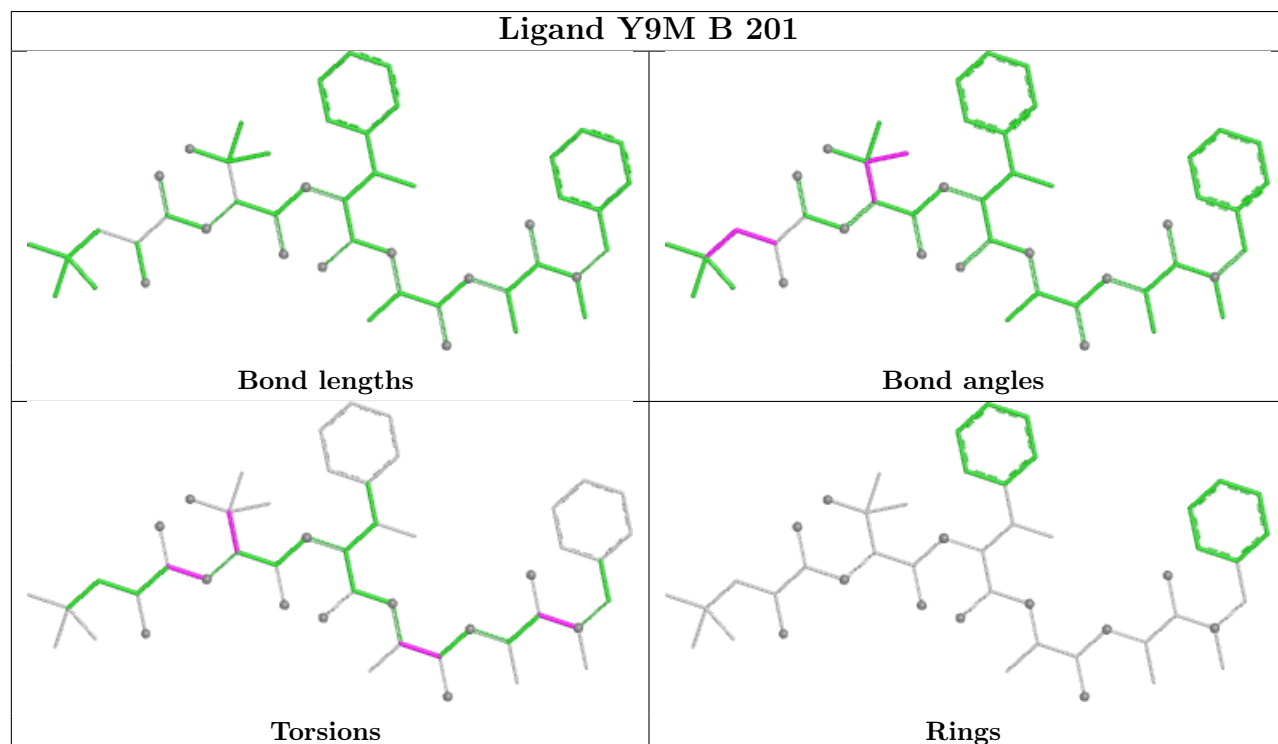
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	201	Y9M	3	0
2	C	201	Y9M	3	0
2	A	201	Y9M	1	0
2	D	201	Y9M	1	0
2	E	201	Y9M	1	0

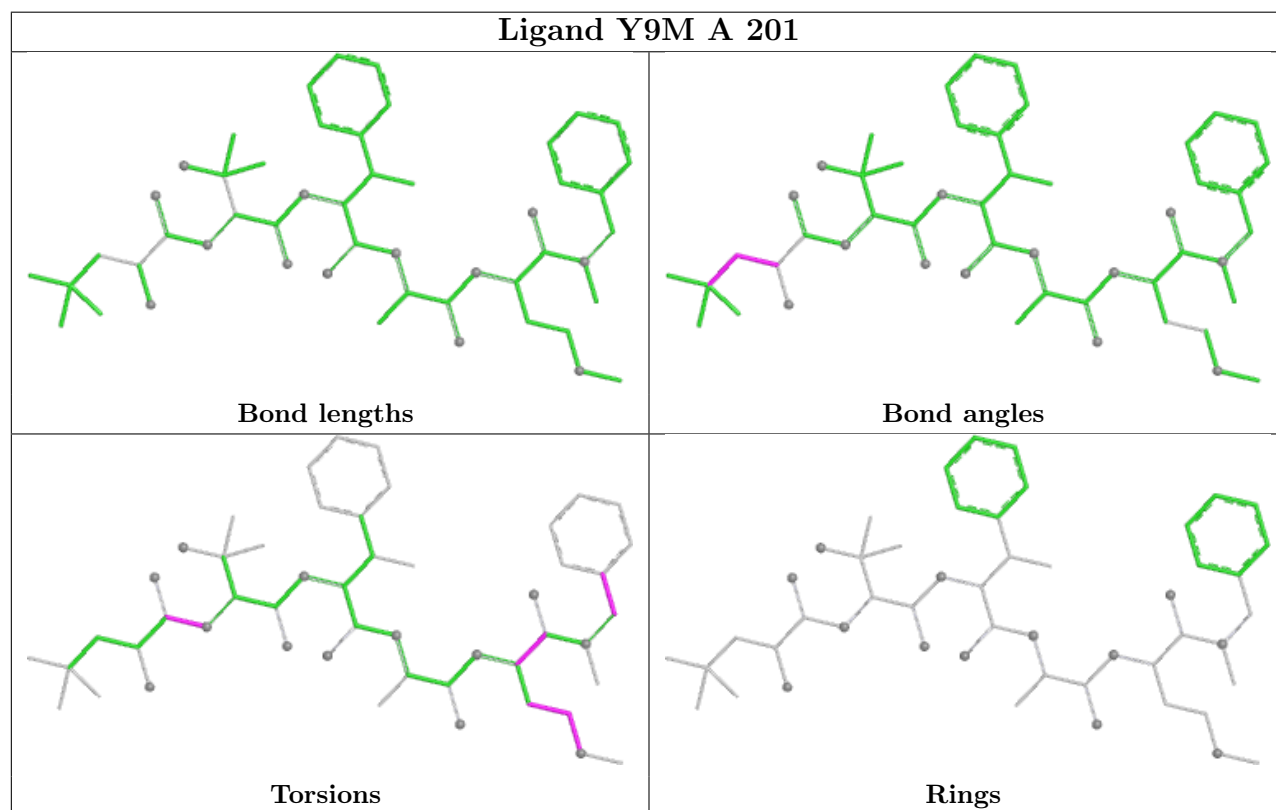
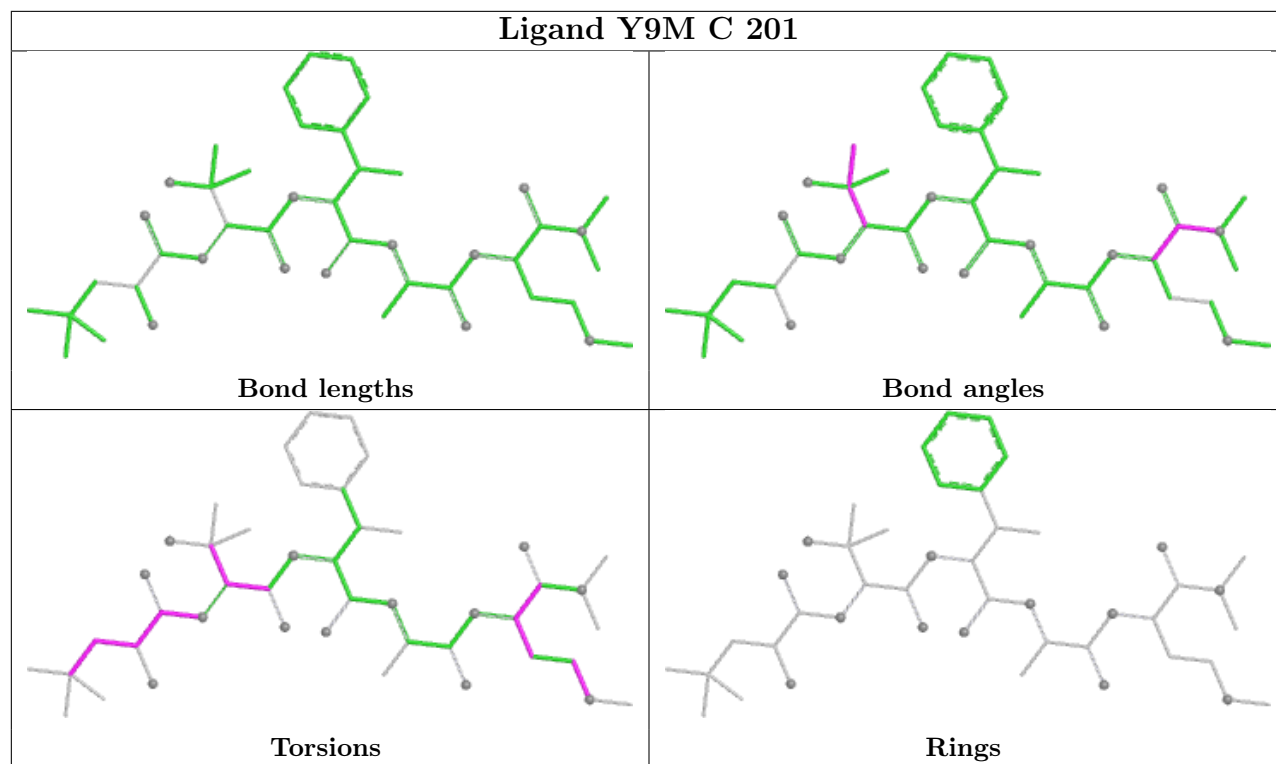
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

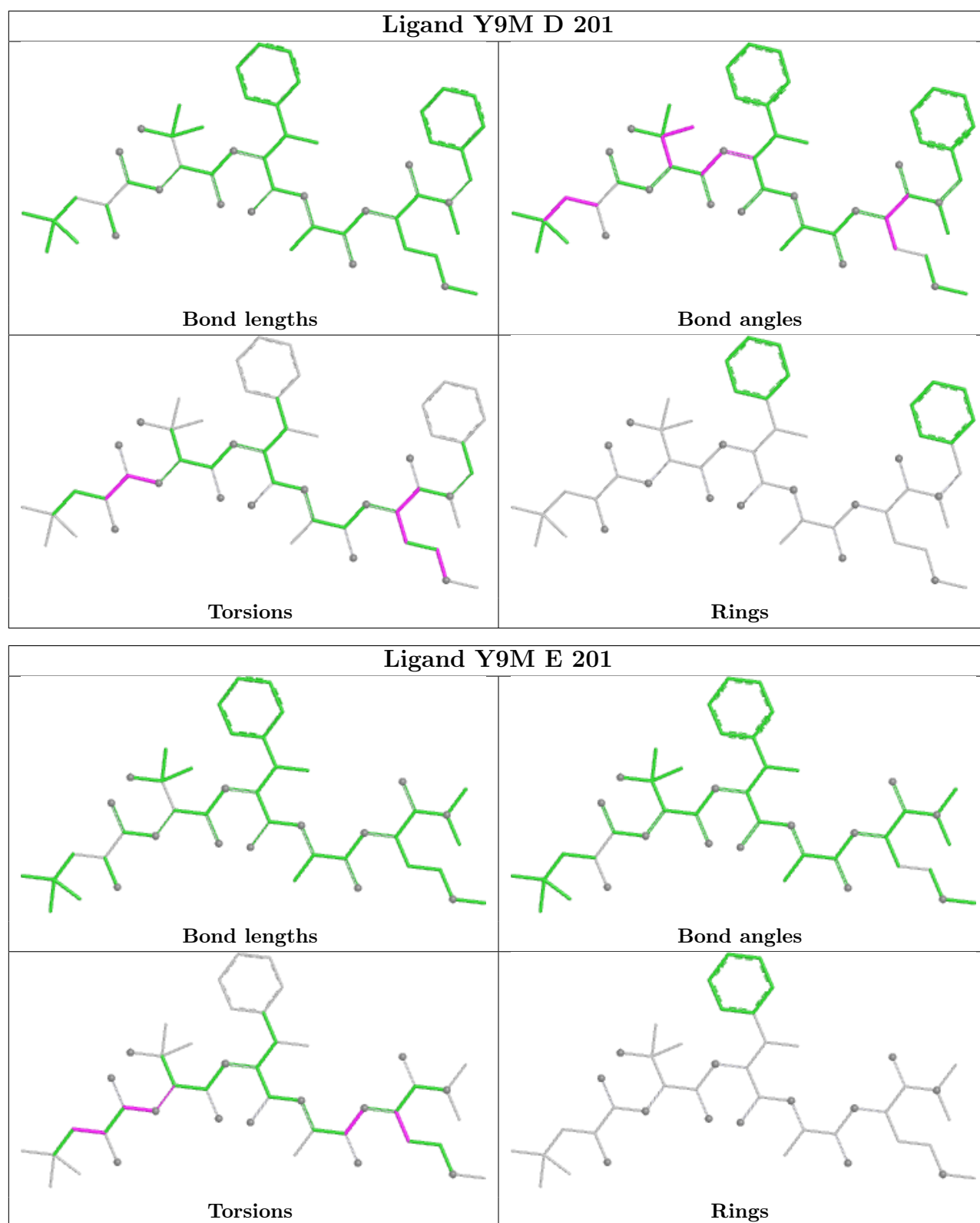
Ligand Y9M F 201



Ligand Y9M B 201







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

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	126/139 (90%)	1.31	23 (18%) 3 3	21, 38, 58, 70	1 (0%)
1	B	133/139 (95%)	0.98	19 (14%) 6 5	18, 34, 50, 63	1 (0%)
1	C	127/139 (91%)	1.18	20 (15%) 5 4	19, 36, 58, 74	2 (1%)
1	D	128/139 (92%)	1.04	12 (9%) 14 12	19, 34, 54, 59	0
1	E	134/139 (96%)	1.24	27 (20%) 3 2	20, 40, 60, 69	2 (1%)
1	F	128/139 (92%)	1.46	31 (24%) 2 1	21, 39, 63, 71	2 (1%)
All	All	776/834 (93%)	1.20	132 (17%) 4 3	18, 37, 59, 74	8 (1%)

All (132) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	25	GLN	5.5
1	F	26	ALA	5.3
1	F	86	LEU	5.2
1	A	125	TYR	5.1
1	E	49	GLY	4.7
1	F	69	ASN	4.6
1	F	42[A]	HIS	4.4
1	A	25	GLN	4.1
1	C	63	ARG	4.0
1	D	81	ASP	3.9
1	F	44	GLY	3.9
1	F	67	SER	3.8
1	E	65	VAL	3.8
1	C	120	VAL	3.7
1	A	19	TYR	3.7
1	A	86	LEU	3.7
1	D	54	ALA	3.6
1	F	19	TYR	3.6
1	E	26	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	F	113	VAL	3.5
1	A	20	TYR	3.4
1	C	29	MET	3.4
1	E	20	TYR	3.4
1	F	22	LEU	3.4
1	E	48	ASN	3.4
1	C	121	ALA	3.3
1	C	73	CYS	3.3
1	E	35	GLY	3.3
1	E	33	PHE	3.3
1	B	120	VAL	3.2
1	A	6	PRO	3.2
1	A	94	VAL	3.2
1	F	29	MET	3.1
1	E	29	MET	3.0
1	E	54	ALA	3.0
1	B	44	GLY	3.0
1	E	50	LYS	2.9
1	A	69	ASN	2.9
1	E	90	VAL	2.9
1	F	23	LEU	2.9
1	A	137	VAL	2.9
1	F	89	GLY	2.8
1	C	122	ASN	2.8
1	A	9	LEU	2.8
1	C	55	VAL	2.8
1	F	2	VAL	2.8
1	A	73	CYS	2.8
1	F	10	LEU	2.8
1	F	28	ASP	2.8
1	F	6	PRO	2.7
1	E	47	SER	2.7
1	F	52	ALA	2.7
1	F	68	GLN	2.7
1	A	26	ALA	2.7
1	C	86	LEU	2.7
1	A	52	ALA	2.7
1	B	39	SER	2.6
1	B	122	ASN	2.6
1	D	72	ASN	2.6
1	B	121	ALA	2.6
1	C	15	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	F	83	HIS	2.6
1	A	34	TYR	2.6
1	B	130	ILE	2.6
1	F	61	ILE	2.6
1	D	19	TYR	2.5
1	E	71	THR	2.5
1	A	85	THR	2.5
1	A	67	SER	2.5
1	C	13	ARG	2.5
1	A	37	ASN	2.5
1	C	130	ILE	2.5
1	B	14	GLU	2.5
1	C	57	GLY	2.5
1	F	53	ASP	2.4
1	C	35	GLY	2.4
1	C	113	VAL	2.4
1	E	4	GLU	2.4
1	D	118	GLY	2.4
1	E	89	GLY	2.4
1	D	9	LEU	2.3
1	B	111	THR	2.3
1	F	66	MET	2.3
1	C	53	ASP	2.3
1	C	33	PHE	2.3
1	B	38	SER	2.3
1	B	88	ASP	2.3
1	F	27	PRO	2.3
1	D	106	ARG	2.3
1	F	93	GLN	2.3
1	E	102	ASN	2.3
1	F	57	GLY	2.3
1	B	113	VAL	2.3
1	E	121	ALA	2.2
1	A	40	TYR	2.2
1	D	40	TYR	2.2
1	A	8	PRO	2.2
1	A	119	SER	2.2
1	E	92	VAL	2.2
1	A	61	ILE	2.2
1	D	69	ASN	2.2
1	E	100	ASN	2.2
1	D	43	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	11	VAL	2.2
1	A	138	PHE	2.2
1	C	37	ASN	2.1
1	E	37	ASN	2.1
1	B	20	TYR	2.1
1	D	92	VAL	2.1
1	F	74	HIS	2.1
1	E	70	PHE	2.1
1	F	33	PHE	2.1
1	B	56	TYR	2.1
1	B	15	PHE	2.1
1	E	73	CYS	2.1
1	F	72	ASN	2.1
1	B	82	ALA	2.1
1	B	58	GLN	2.1
1	A	24	ASN	2.1
1	B	43	GLY	2.1
1	E	74	HIS	2.1
1	E	137	VAL	2.1
1	E	36	LYS	2.0
1	F	103	GLN	2.0
1	E	31	HIS	2.0
1	E	19	TYR	2.0
1	C	38	SER	2.0
1	D	110	GLN	2.0
1	F	105	LEU	2.0
1	B	12	GLY	2.0
1	B	57	GLY	2.0
1	C	102	ASN	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 ⓘ

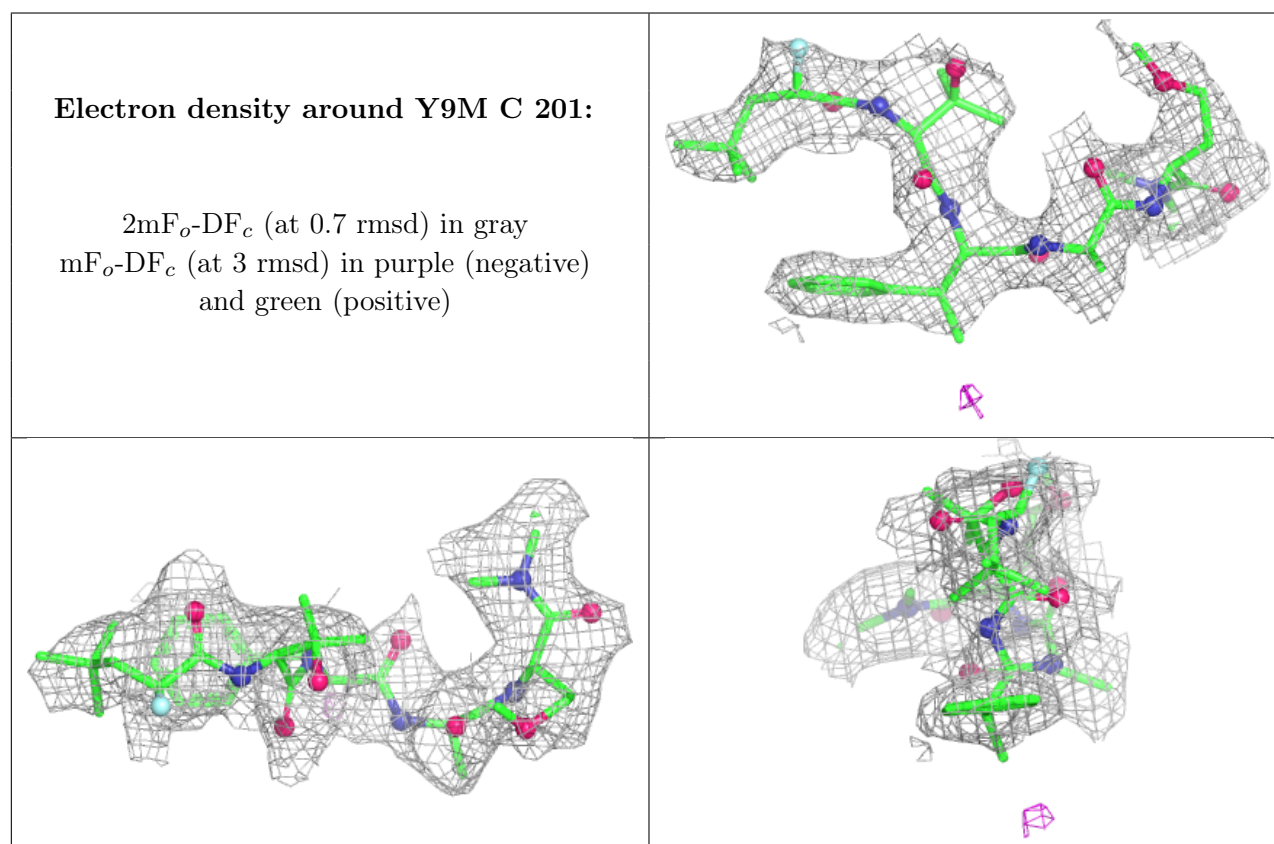
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

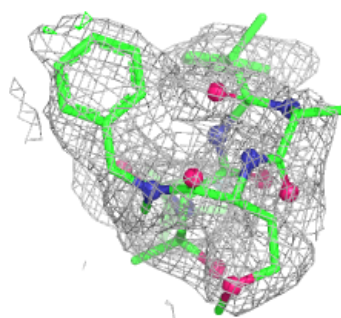
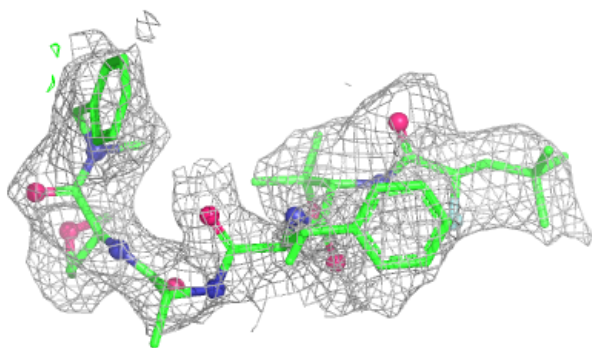
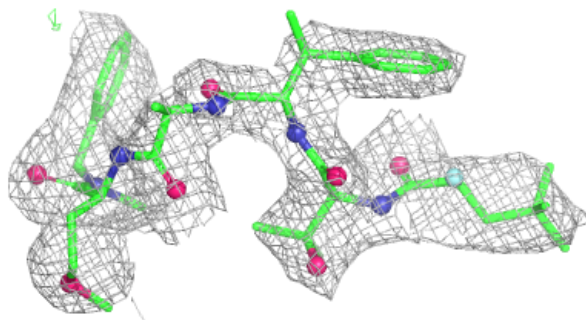
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	Y9M	C	201	45/51	0.78	0.15	44,52,57,59	0
2	Y9M	D	201	51/51	0.79	0.16	34,41,52,54	0
2	Y9M	A	201	51/51	0.80	0.17	35,47,57,59	0
2	Y9M	B	201	48/51	0.80	0.18	39,53,60,62	0
2	Y9M	E	201	45/51	0.82	0.15	34,43,59,69	0
2	Y9M	F	201	51/51	0.84	0.15	30,39,54,57	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.

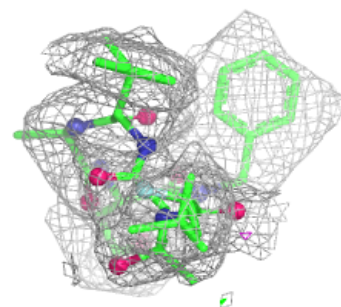
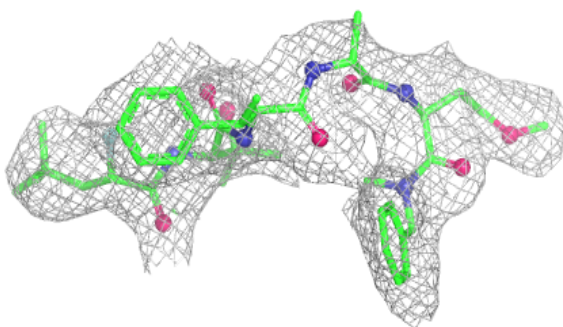
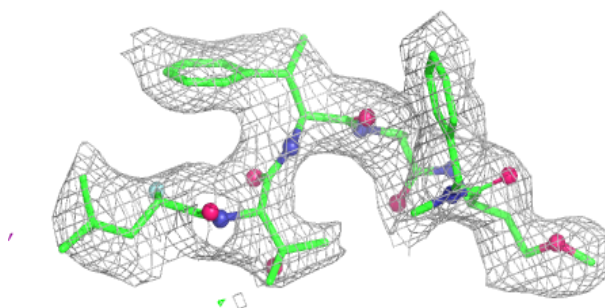


Electron density around Y9M D 201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

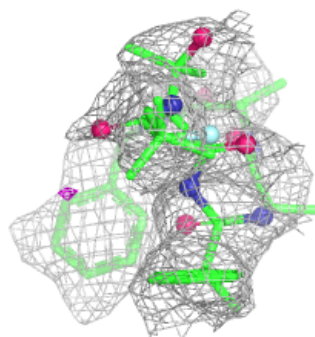
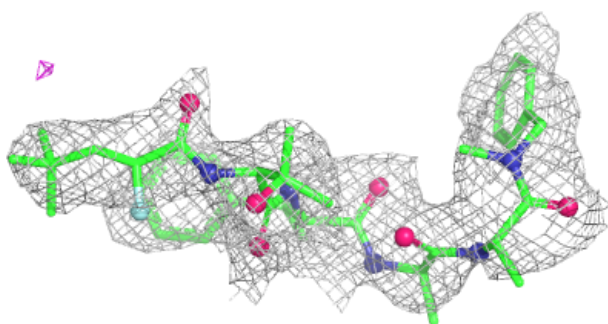
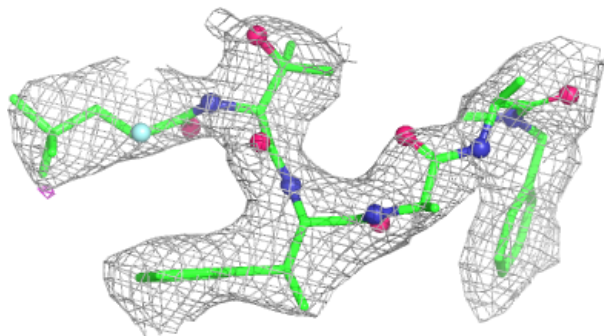
**Electron density around Y9M A 201:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

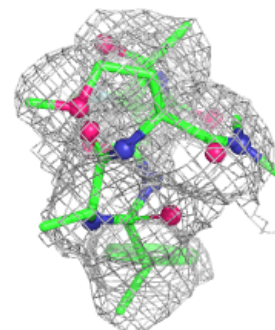
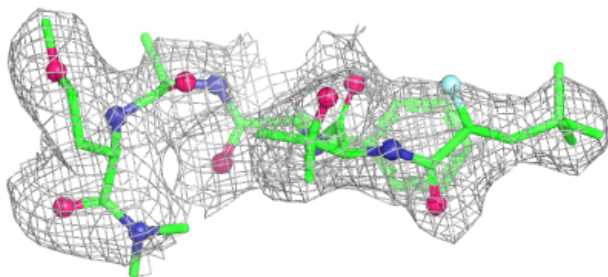
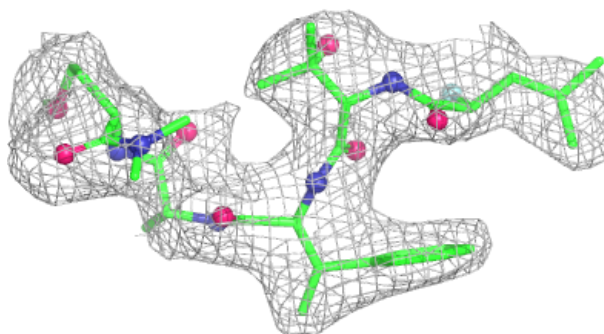


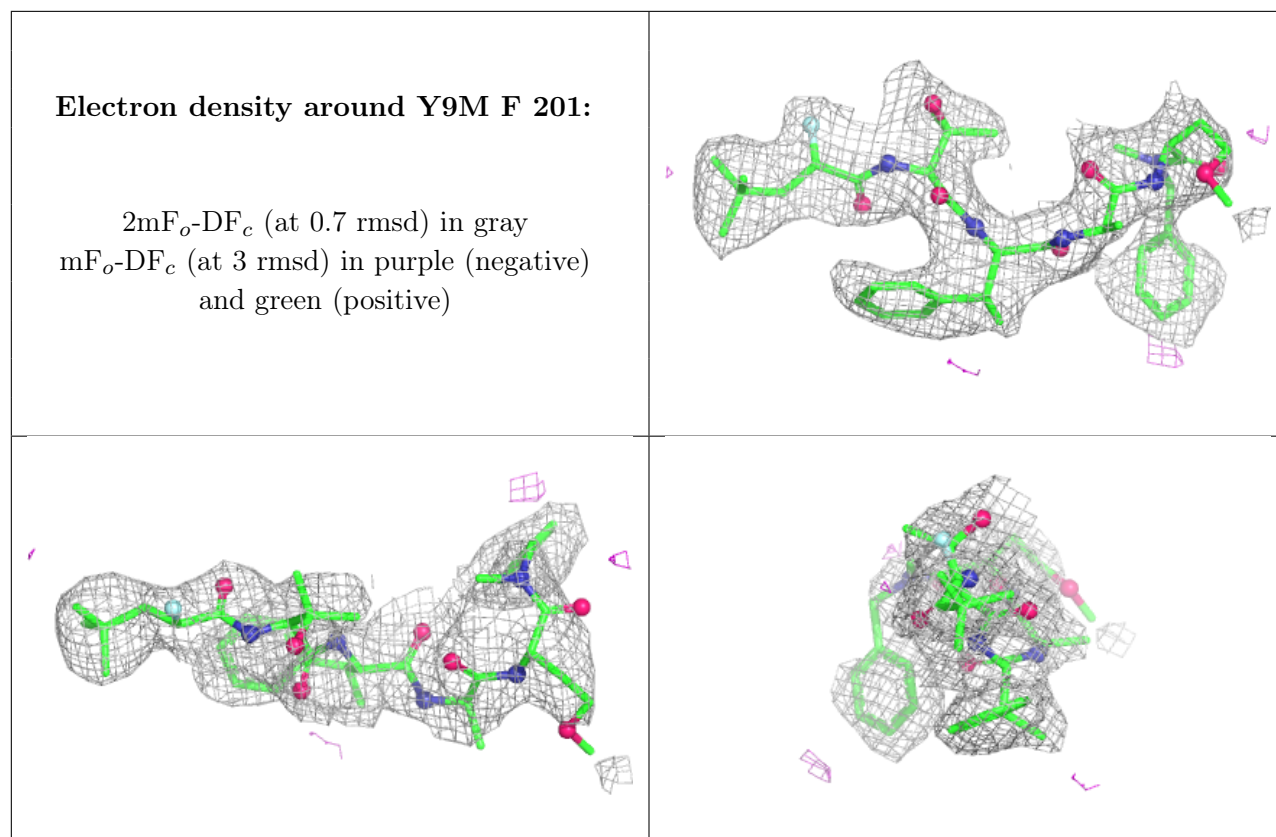
Electron density around Y9M B 201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around Y9M E 201:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.