



Full wwPDB/EMDatabank EM Map/Model Validation Report ⓘ

Jan 22, 2018 – 09:09 PM EST

PDB ID : 6B5V
EMDB ID: : EMD-7058
Title : Structure of TRPV5 in complex with econazole
Authors : Hughes, T.E.T.; Lodowski, D.T.; Huynh, K.W.; Yazici, A.; del Rosario, J.; Kapoor, A.; Basak, S.; Samanta, A.; Chakrapani, S.; Zhou, Z.H.; Filizola, M.; Rohacs, T.; Han, S.; Moiseenkova-Bell, V.Y.
Deposited on : 2017-09-29
Resolution : 4.80 Å(reported)

This is a Full wwPDB/EMDatabank EM Map/Model Validation Report
for a publicly released PDB/EMDB entry.

We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030736

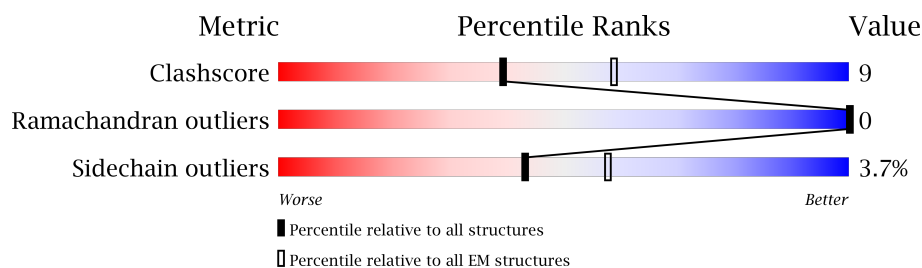
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.80 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)	EM structures (#Entries)
Clashscore	125131	1336
Ramachandran outliers	121729	1120
Sidechain outliers	121581	1026

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	730	66% 14% • 19%
1	B	730	66% 14% • 19%
1	C	730	66% 14% • 19%
1	D	730	66% 14% • 19%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ECL	A	1001	-	-	X	-
2	ECL	B	1001	-	-	X	-
2	ECL	C	1001	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ECL	D	1001	-	-	X	-

2 Entry composition [i](#)

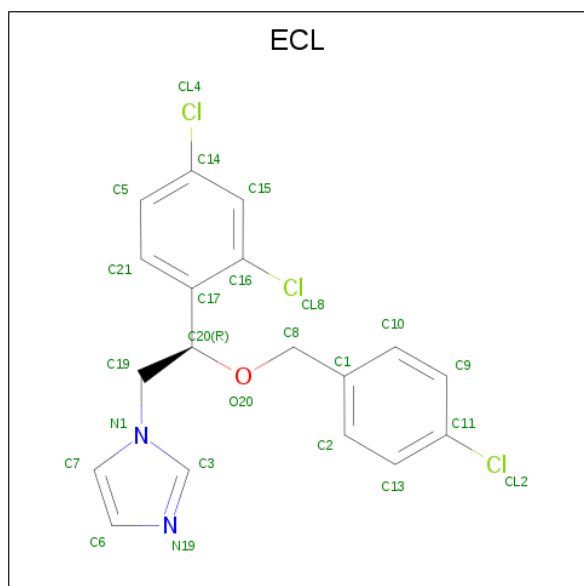
There are 3 unique types of molecules in this entry. The entry contains 18993 atoms, of which 60 are hydrogens and 0 are deuteriums.

In the tables below, 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 Transient receptor potential cation channel subfamily V member 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	589	Total	C	N	O	S	0	0
			4709	3051	795	828	35		
1	B	589	Total	C	N	O	S	0	0
			4709	3051	795	828	35		
1	D	589	Total	C	N	O	S	0	0
			4709	3051	795	828	35		
1	C	589	Total	C	N	O	S	0	0
			4709	3051	795	828	35		

- Molecule 2 is 1-[(2R)-2-[(4-chlorobenzyl)oxy]-2-(2,4-dichlorophenyl)ethyl]-1H-imidazole (three-letter code: ECL) (formula: C₁₈H₁₅Cl₃N₂O).



Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	Cl	H	N	O
			39	18	3	15	2	1
2	B	1	Total	C	Cl	H	N	O
			39	18	3	15	2	1

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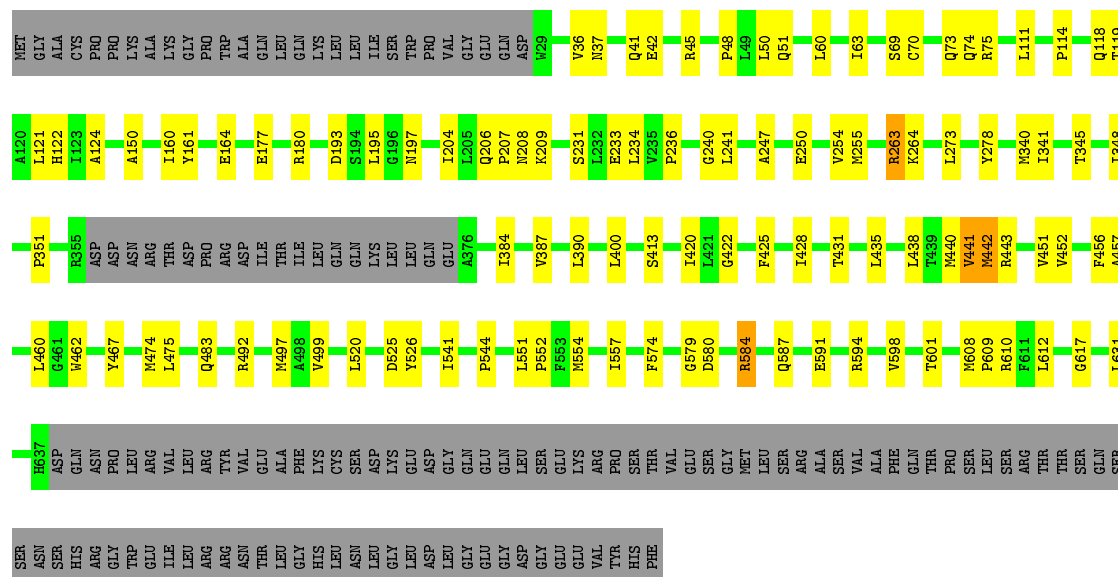
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Mol	Chain	Residues	Atoms						AltConf
2	D	1	Total	C	Cl	H	N	O	0
			39	18	3	15	2	1	
2	C	1	Total	C	Cl	H	N	O	0
			39	18	3	15	2	1	

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
3	A	1	Total	Ca	0
			1	1	

- Chain D:  66% 14% • 19%



4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C4	Depositor
Number of particles used	50566	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	2	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	45455	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	0.31	0/4821	0.66	0/6546
1	B	0.31	0/4821	0.66	0/6546
1	C	0.31	0/4821	0.66	0/6546
1	D	0.31	0/4821	0.66	0/6546
All	All	0.31	0/19284	0.66	0/26184

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4709	0	4741	82	0
1	B	4709	0	4741	85	0
1	C	4709	0	4741	81	0
1	D	4709	0	4741	82	0
2	A	24	15	15	8	0
2	B	24	15	15	8	0
2	C	24	15	15	8	0
2	D	24	15	15	8	0
3	A	1	0	0	0	0
All	All	18933	60	19024	338	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 9.

All (338) 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:263:ARG:HD3	1:B:263:ARG:C	1.59	1.20
1:A:263:ARG:C	1:A:263:ARG:HD3	1.59	1.20
1:C:263:ARG:C	1:C:263:ARG:HD3	1.59	1.19
1:D:263:ARG:C	1:D:263:ARG:HD3	1.59	1.13
1:A:435:LEU:HD22	1:A:457:ALA:CB	1.94	0.97
1:B:435:LEU:HD22	1:B:457:ALA:CB	1.94	0.97
1:C:435:LEU:HD22	1:C:457:ALA:CB	1.94	0.96
1:D:435:LEU:HD22	1:D:457:ALA:CB	1.94	0.96
1:A:45:ARG:HG2	1:A:45:ARG:HH11	1.34	0.92
1:D:45:ARG:HG2	1:D:45:ARG:HH11	1.34	0.92
1:C:45:ARG:HH11	1:C:45:ARG:HG2	1.34	0.92
1:B:45:ARG:HG2	1:B:45:ARG:HH11	1.34	0.91
1:C:435:LEU:HD22	1:C:457:ALA:HB3	1.53	0.91
1:D:435:LEU:HD22	1:D:457:ALA:HB3	1.53	0.89
1:A:435:LEU:HD22	1:A:457:ALA:HB3	1.53	0.88
1:C:263:ARG:CD	1:C:263:ARG:C	2.41	0.88
1:B:435:LEU:HD22	1:B:457:ALA:HB3	1.53	0.87
1:D:263:ARG:C	1:D:263:ARG:CD	2.41	0.84
1:B:263:ARG:C	1:B:263:ARG:CD	2.41	0.83
1:D:610:ARG:HE	1:D:610:ARG:HA	1.44	0.82
1:A:610:ARG:HA	1:A:610:ARG:HE	1.44	0.82
1:B:610:ARG:HE	1:B:610:ARG:HA	1.44	0.81
1:C:610:ARG:HE	1:C:610:ARG:HA	1.44	0.80
1:B:177:GLU:O	1:B:180:ARG:HB3	1.83	0.79
1:D:263:ARG:HD3	1:D:263:ARG:O	1.82	0.79
1:B:263:ARG:HD3	1:B:263:ARG:O	1.82	0.79
1:D:177:GLU:O	1:D:180:ARG:HB3	1.83	0.78
1:A:177:GLU:O	1:A:180:ARG:HB3	1.83	0.78
1:A:263:ARG:O	1:A:263:ARG:HD3	1.82	0.78
1:C:263:ARG:O	1:C:263:ARG:HD3	1.82	0.78
1:C:177:GLU:O	1:C:180:ARG:HB3	1.83	0.77
1:D:610:ARG:NE	1:D:610:ARG:HA	2.01	0.73
1:A:263:ARG:CD	1:A:263:ARG:C	2.41	0.72
1:C:435:LEU:O	1:C:435:LEU:HD12	1.90	0.72
1:A:610:ARG:HA	1:A:610:ARG:NE	2.01	0.71
1:B:263:ARG:HD3	1:B:264:LYS:N	2.05	0.71
1:D:263:ARG:HD3	1:D:264:LYS:N	2.05	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:435:LEU:O	1:D:435:LEU:HD12	1.90	0.71
1:A:263:ARG:HD3	1:A:264:LYS:N	2.05	0.71
1:C:263:ARG:HD3	1:C:264:LYS:N	2.05	0.71
1:A:435:LEU:O	1:A:435:LEU:HD12	1.90	0.70
1:B:435:LEU:HD12	1:B:435:LEU:O	1.90	0.70
1:B:610:ARG:HA	1:B:610:ARG:NE	2.01	0.70
1:C:610:ARG:NE	1:C:610:ARG:HA	2.01	0.69
1:A:263:ARG:CD	1:A:263:ARG:O	2.43	0.67
1:B:263:ARG:O	1:B:263:ARG:CD	2.43	0.66
1:C:263:ARG:CD	1:C:263:ARG:O	2.43	0.66
1:D:341:ILE:O	1:D:345:THR:HG23	1.96	0.66
1:A:435:LEU:CD2	1:A:457:ALA:HB3	2.26	0.66
1:D:263:ARG:CD	1:D:263:ARG:O	2.43	0.66
1:A:341:ILE:O	1:A:345:THR:HG23	1.95	0.66
1:B:435:LEU:CD2	1:B:457:ALA:HB3	2.25	0.65
1:B:341:ILE:O	1:B:345:THR:HG23	1.95	0.65
1:C:341:ILE:O	1:C:345:THR:HG23	1.95	0.65
1:B:45:ARG:NH1	1:B:45:ARG:HG2	2.10	0.65
1:D:435:LEU:CD2	1:D:457:ALA:HB3	2.26	0.64
1:C:435:LEU:CD2	1:C:457:ALA:HB3	2.26	0.64
2:C:1001:ECL:H7	2:C:1001:ECL:H21	1.81	0.63
2:A:1001:ECL:H21	2:A:1001:ECL:H7	1.81	0.63
2:B:1001:ECL:H7	2:B:1001:ECL:H21	1.80	0.63
1:A:442:MET:HG3	1:A:443:ARG:HG3	1.81	0.63
1:C:442:MET:HG3	1:C:443:ARG:HG3	1.81	0.62
1:D:442:MET:HG3	1:D:443:ARG:HG3	1.81	0.62
1:B:442:MET:HG3	1:B:443:ARG:HG3	1.81	0.62
1:A:456:PHE:HB2	1:C:554:MET:HE1	1.82	0.62
2:D:1001:ECL:H7	2:D:1001:ECL:H21	1.80	0.62
1:A:617:GLY:HA2	1:A:631:LEU:HD23	1.82	0.62
1:B:45:ARG:CG	1:B:45:ARG:HH11	2.11	0.62
1:C:617:GLY:HA2	1:C:631:LEU:HD23	1.82	0.61
1:D:345:THR:O	1:D:348:ILE:HG12	2.00	0.61
1:A:345:THR:O	1:A:348:ILE:HG12	2.01	0.61
1:A:45:ARG:CG	1:A:45:ARG:HH11	2.11	0.61
1:B:345:THR:O	1:B:348:ILE:HG12	2.01	0.61
1:B:617:GLY:HA2	1:B:631:LEU:HD23	1.82	0.61
1:D:617:GLY:HA2	1:D:631:LEU:HD23	1.82	0.61
1:C:45:ARG:NH1	1:C:45:ARG:HG2	2.10	0.61
1:B:579:GLY:O	1:B:584:ARG:HG3	2.01	0.60
1:C:579:GLY:O	1:C:584:ARG:HG3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:579:GLY:O	1:D:584:ARG:HG3	2.01	0.60
1:B:438:LEU:HA	1:B:441:VAL:HG13	1.83	0.60
1:A:579:GLY:O	1:A:584:ARG:HG3	2.01	0.59
1:A:438:LEU:HA	1:A:441:VAL:HG13	1.83	0.59
1:C:345:THR:O	1:C:348:ILE:HG12	2.01	0.59
1:D:438:LEU:HA	1:D:441:VAL:HG13	1.83	0.59
1:C:438:LEU:HA	1:C:441:VAL:HG13	1.83	0.58
1:D:612:LEU:O	1:D:612:LEU:HG	2.04	0.58
1:D:580:ASP:HA	1:D:584:ARG:HB2	1.85	0.58
1:B:42:GLU:HA	1:B:45:ARG:HG3	1.85	0.58
1:A:612:LEU:O	1:A:612:LEU:HG	2.04	0.58
1:B:612:LEU:HG	1:B:612:LEU:O	2.04	0.58
1:C:612:LEU:HG	1:C:612:LEU:O	2.04	0.58
1:B:580:ASP:HA	1:B:584:ARG:HB2	1.85	0.57
1:C:580:ASP:HA	1:C:584:ARG:HB2	1.85	0.57
1:C:45:ARG:CG	1:C:45:ARG:HH11	2.11	0.57
2:A:1001:ECL:CL8	2:A:1001:ECL:C8	2.89	0.57
1:C:42:GLU:HA	1:C:45:ARG:CG	2.34	0.57
1:C:42:GLU:HB3	1:C:45:ARG:HD3	1.86	0.57
2:B:1001:ECL:C8	2:B:1001:ECL:CL8	2.89	0.57
1:A:42:GLU:HB3	1:A:45:ARG:HD3	1.86	0.57
1:C:42:GLU:HA	1:C:45:ARG:HG3	1.85	0.57
2:D:1001:ECL:CL8	2:D:1001:ECL:C8	2.89	0.57
1:A:42:GLU:HA	1:A:45:ARG:HG3	1.85	0.57
1:B:42:GLU:HA	1:B:45:ARG:CG	2.34	0.57
2:C:1001:ECL:CL8	2:C:1001:ECL:C8	2.89	0.57
1:B:554:MET:HE3	1:B:557:ILE:HD11	1.87	0.57
1:A:580:ASP:HA	1:A:584:ARG:HB2	1.85	0.56
1:A:42:GLU:HA	1:A:45:ARG:CG	2.34	0.56
1:D:42:GLU:HA	1:D:45:ARG:HG3	1.85	0.56
1:B:161:TYR:OH	1:B:197:ASN:ND2	2.39	0.56
1:D:42:GLU:HA	1:D:45:ARG:CG	2.34	0.56
1:A:161:TYR:OH	1:A:197:ASN:ND2	2.39	0.56
1:D:161:TYR:OH	1:D:197:ASN:ND2	2.39	0.56
1:B:42:GLU:HB3	1:B:45:ARG:HD3	1.86	0.55
1:D:42:GLU:HB3	1:D:45:ARG:HD3	1.86	0.55
1:A:45:ARG:NH1	1:A:45:ARG:HG2	2.10	0.55
1:A:384:ILE:HA	1:A:387:VAL:HG22	1.89	0.55
1:B:384:ILE:HA	1:B:387:VAL:HG22	1.89	0.55
1:D:554:MET:HE1	1:C:456:PHE:HB2	1.88	0.55
1:D:520:LEU:HD11	1:D:544:PRO:HB2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:384:ILE:HA	1:C:387:VAL:HG22	1.89	0.55
1:D:384:ILE:HA	1:D:387:VAL:HG22	1.89	0.55
1:A:420:ILE:HG23	1:A:483:GLN:HE22	1.72	0.55
1:C:161:TYR:OH	1:C:197:ASN:ND2	2.39	0.54
1:A:520:LEU:HD11	1:A:544:PRO:HB2	1.89	0.54
1:B:520:LEU:HD11	1:B:544:PRO:HB2	1.89	0.54
1:D:45:ARG:HG2	1:D:45:ARG:NH1	2.10	0.54
1:D:420:ILE:HG23	1:D:483:GLN:HE22	1.72	0.54
1:C:520:LEU:HD11	1:C:544:PRO:HB2	1.89	0.54
1:B:42:GLU:CB	1:B:45:ARG:HD3	2.39	0.53
1:B:420:ILE:HG23	1:B:483:GLN:HE22	1.72	0.53
1:A:42:GLU:CB	1:A:45:ARG:HD3	2.39	0.53
1:C:420:ILE:HG23	1:C:483:GLN:HE22	1.72	0.53
1:D:42:GLU:CB	1:D:45:ARG:HD3	2.39	0.53
1:B:554:MET:HE1	1:D:456:PHE:HB2	1.91	0.53
1:A:554:MET:HE2	1:B:456:PHE:HB2	1.91	0.52
1:D:45:ARG:CG	1:D:45:ARG:HH11	2.11	0.52
1:C:42:GLU:CB	1:C:45:ARG:HD3	2.39	0.52
1:B:551:LEU:HD22	1:B:552:PRO:HD2	1.92	0.52
1:A:551:LEU:HD22	1:A:552:PRO:HD2	1.92	0.51
2:A:1001:ECL:H21	2:A:1001:ECL:C7	2.40	0.51
1:D:499:VAL:O	1:C:462:TRP:NE1	2.42	0.51
1:C:45:ARG:CG	1:C:45:ARG:NH1	2.72	0.51
2:B:1001:ECL:C7	2:B:1001:ECL:H21	2.40	0.51
1:B:499:VAL:O	1:D:462:TRP:NE1	2.44	0.51
1:D:580:ASP:O	1:D:584:ARG:HB2	2.11	0.51
1:B:263:ARG:O	1:B:263:ARG:CG	2.60	0.50
2:C:1001:ECL:C7	2:C:1001:ECL:H21	2.40	0.50
2:D:1001:ECL:C7	2:D:1001:ECL:H21	2.40	0.50
1:C:263:ARG:CG	1:C:263:ARG:O	2.60	0.50
1:C:551:LEU:HD22	1:C:552:PRO:HD2	1.92	0.50
1:D:554:MET:HE3	1:D:557:ILE:HD11	1.93	0.50
1:A:45:ARG:CG	1:A:45:ARG:NH1	2.72	0.50
1:B:580:ASP:O	1:B:584:ARG:HB2	2.12	0.50
1:D:551:LEU:HD22	1:D:552:PRO:HD2	1.92	0.50
1:D:263:ARG:CG	1:D:263:ARG:O	2.60	0.50
1:C:351:PRO:HG3	1:C:451:VAL:HG11	1.94	0.50
1:A:387:VAL:HA	1:A:390:LEU:HB2	1.94	0.50
1:B:387:VAL:HA	1:B:390:LEU:HB2	1.94	0.50
1:C:580:ASP:O	1:C:584:ARG:HB2	2.12	0.50
1:D:351:PRO:HG3	1:D:451:VAL:HG11	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121:LEU:O	1:B:124:ALA:N	2.46	0.49
1:B:37:ASN:O	1:B:41:GLN:NE2	2.43	0.49
1:B:69:SER:O	1:B:73:GLN:NE2	2.45	0.49
1:C:387:VAL:HA	1:C:390:LEU:HB2	1.94	0.49
1:A:263:ARG:O	1:A:263:ARG:CG	2.60	0.49
1:A:69:SER:O	1:A:73:GLN:NE2	2.46	0.49
1:A:351:PRO:HG3	1:A:451:VAL:HG11	1.94	0.49
1:A:610:ARG:HD3	1:A:610:ARG:O	2.13	0.49
1:C:278:TYR:HB2	1:C:631:LEU:HB3	1.94	0.49
1:A:580:ASP:O	1:A:584:ARG:HB2	2.12	0.49
1:A:209:LYS:HG3	1:A:254:VAL:HG11	1.95	0.49
1:C:610:ARG:O	1:C:610:ARG:HD3	2.13	0.49
1:C:69:SER:O	1:C:73:GLN:NE2	2.45	0.49
1:D:69:SER:O	1:D:73:GLN:NE2	2.45	0.49
1:D:278:TYR:HB2	1:D:631:LEU:HB3	1.94	0.49
1:D:387:VAL:HA	1:D:390:LEU:HB2	1.94	0.49
1:B:610:ARG:HD3	1:B:610:ARG:O	2.13	0.49
1:B:278:TYR:HB2	1:B:631:LEU:HB3	1.94	0.48
1:A:435:LEU:CD2	1:A:457:ALA:CB	2.81	0.48
1:C:121:LEU:O	1:C:124:ALA:N	2.46	0.48
1:C:60:LEU:HD23	1:C:63:ILE:HD12	1.96	0.48
1:D:209:LYS:HG3	1:D:254:VAL:HG11	1.95	0.48
1:A:204:ILE:HG22	1:A:255:MET:HG3	1.96	0.48
1:A:278:TYR:HB2	1:A:631:LEU:HB3	1.94	0.48
1:A:60:LEU:HD23	1:A:63:ILE:HD12	1.96	0.48
1:B:204:ILE:HG22	1:B:255:MET:HG3	1.96	0.48
1:B:209:LYS:HG3	1:B:254:VAL:HG11	1.95	0.48
1:B:525:ASP:OD1	1:B:526:TYR:N	2.47	0.48
1:D:610:ARG:O	1:D:610:ARG:HD3	2.13	0.48
1:A:160:ILE:HG21	1:B:273:LEU:HD11	1.96	0.48
1:A:462:TRP:NE1	1:C:499:VAL:O	2.46	0.48
1:D:231:SER:HB3	1:D:234:LEU:HB2	1.96	0.48
1:D:60:LEU:HD23	1:D:63:ILE:HD12	1.96	0.48
1:B:351:PRO:HG3	1:B:451:VAL:HG11	1.94	0.48
1:C:204:ILE:HG22	1:C:255:MET:HG3	1.96	0.48
1:A:121:LEU:O	1:A:124:ALA:N	2.46	0.48
1:B:460:LEU:HD11	2:B:1001:ECL:H13	1.96	0.48
1:C:209:LYS:HG3	1:C:254:VAL:HG11	1.95	0.48
1:D:460:LEU:HD11	2:D:1001:ECL:H13	1.96	0.48
1:C:428:ILE:HA	1:C:431:THR:HG22	1.96	0.47
1:B:60:LEU:HD23	1:B:63:ILE:HD12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:121:LEU:O	1:D:124:ALA:N	2.46	0.47
1:D:428:ILE:HA	1:D:431:THR:HG22	1.96	0.47
1:C:460:LEU:HD11	2:C:1001:ECL:H13	1.96	0.47
1:D:204:ILE:HG22	1:D:255:MET:HG3	1.96	0.47
1:A:460:LEU:HD11	2:A:1001:ECL:H13	1.96	0.47
1:C:37:ASN:O	1:C:41:GLN:NE2	2.43	0.47
1:A:50:LEU:HD13	1:A:74:GLN:HB3	1.97	0.47
1:B:231:SER:HB3	1:B:234:LEU:HB2	1.96	0.47
1:D:50:LEU:HD13	1:D:74:GLN:HB3	1.97	0.47
1:A:231:SER:HB3	1:A:234:LEU:HB2	1.96	0.47
1:B:425:PHE:CD2	1:B:467:TYR:HD1	2.33	0.47
1:A:428:ILE:HA	1:A:431:THR:HG22	1.96	0.47
1:B:428:ILE:HA	1:B:431:THR:HG22	1.96	0.47
1:B:237:ASN:OD1	1:B:241:LEU:N	2.47	0.47
1:C:50:LEU:HD13	1:C:74:GLN:HB3	1.97	0.47
1:D:525:ASP:OD1	1:D:526:TYR:N	2.47	0.47
1:C:164:GLU:HB2	1:C:193:ASP:HB3	1.98	0.46
1:C:425:PHE:CD2	1:C:467:TYR:HD1	2.33	0.46
1:C:231:SER:HB3	1:C:234:LEU:HB2	1.96	0.46
1:A:164:GLU:HB2	1:A:193:ASP:HB3	1.98	0.46
1:B:50:LEU:HD13	1:B:74:GLN:HB3	1.97	0.46
1:C:236:PRO:HB2	1:C:240:GLY:HA2	1.98	0.46
1:D:164:GLU:HB2	1:D:193:ASP:HB3	1.98	0.46
1:A:425:PHE:CD2	1:A:467:TYR:HD1	2.33	0.46
1:C:525:ASP:OD1	1:C:526:TYR:N	2.47	0.46
1:A:598:VAL:HA	1:A:601:THR:HG22	1.98	0.46
1:D:598:VAL:HA	1:D:601:THR:HG22	1.98	0.46
2:D:1001:ECL:H19	2:D:1001:ECL:H2	1.97	0.46
1:B:164:GLU:HB2	1:B:193:ASP:HB3	1.97	0.46
1:B:428:ILE:HG13	1:B:428:ILE:H	1.61	0.46
1:D:425:PHE:CD2	1:D:467:TYR:HD1	2.33	0.46
1:C:598:VAL:HA	1:C:601:THR:HG22	1.98	0.46
1:D:236:PRO:HB2	1:D:240:GLY:HA2	1.98	0.46
1:A:525:ASP:OD1	1:A:526:TYR:N	2.47	0.45
2:B:1001:ECL:H2	2:B:1001:ECL:H19	1.97	0.45
1:B:45:ARG:NH1	1:B:45:ARG:CG	2.72	0.45
1:B:598:VAL:HA	1:B:601:THR:HG22	1.98	0.45
1:D:237:ASN:OD1	1:D:241:LEU:N	2.47	0.45
1:A:236:PRO:HB2	1:A:240:GLY:HA2	1.98	0.45
2:C:1001:ECL:H19	2:C:1001:ECL:H2	1.97	0.45
2:A:1001:ECL:H2	2:A:1001:ECL:H19	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:233:GLU:HG2	1:C:234:LEU:HD12	1.99	0.45
1:A:37:ASN:HD22	1:B:632:ARG:NH1	2.15	0.45
1:A:233:GLU:HG2	1:A:234:LEU:HD12	1.99	0.45
1:A:88:LEU:HD21	1:B:268:TRP:HB2	1.99	0.45
1:C:428:ILE:HG13	1:C:428:ILE:H	1.62	0.45
1:D:542:ASP:HB2	1:C:541:ILE:HG22	1.99	0.45
1:B:236:PRO:HB2	1:B:240:GLY:HA2	1.98	0.45
1:A:37:ASN:O	1:A:41:GLN:NE2	2.43	0.45
1:A:428:ILE:HG13	1:A:428:ILE:H	1.62	0.44
1:B:233:GLU:HG2	1:B:234:LEU:HD12	1.98	0.44
1:C:193:ASP:OD1	1:C:197:ASN:N	2.50	0.44
1:A:591:GLU:HA	1:A:594:ARG:HG2	2.00	0.44
1:C:591:GLU:HA	1:C:594:ARG:HG2	2.00	0.44
1:D:193:ASP:OD1	1:D:197:ASN:N	2.50	0.44
1:B:193:ASP:OD1	1:B:197:ASN:N	2.50	0.44
1:B:591:GLU:HA	1:B:594:ARG:HG2	2.00	0.44
1:D:591:GLU:HA	1:D:594:ARG:HG2	2.00	0.44
1:A:193:ASP:OD1	1:A:197:ASN:N	2.50	0.44
1:A:206:GLN:HA	1:A:207:PRO:HD3	1.84	0.44
1:C:608:MET:HA	1:C:609:PRO:HD3	1.81	0.44
1:D:233:GLU:HG2	1:D:234:LEU:HD12	1.98	0.44
2:A:1001:ECL:CL8	2:A:1001:ECL:H8A	2.56	0.43
1:C:206:GLN:HA	1:C:207:PRO:HD3	1.84	0.43
1:B:202:ILE:H	1:B:202:ILE:HG13	1.71	0.43
1:D:37:ASN:O	1:D:41:GLN:NE2	2.43	0.43
1:B:206:GLN:HE22	1:B:208:ASN:HB2	1.84	0.43
1:B:301:LYS:HZ3	1:B:303:GLU:HB3	1.83	0.43
1:A:206:GLN:HE22	1:A:208:ASN:HB2	1.84	0.43
2:B:1001:ECL:O20	2:B:1001:ECL:CL8	2.74	0.43
2:B:1001:ECL:H8A	2:B:1001:ECL:CL8	2.56	0.43
1:A:75:ARG:HH22	1:A:111:LEU:HD23	1.84	0.43
1:B:48:PRO:HA	1:B:51:GLN:HE22	1.84	0.43
1:C:263:ARG:NH2	1:C:278:TYR:HA	2.34	0.43
1:C:75:ARG:HH22	1:C:111:LEU:HD23	1.84	0.43
2:C:1001:ECL:CL8	2:C:1001:ECL:H8A	2.56	0.43
1:D:48:PRO:HA	1:D:51:GLN:HE22	1.84	0.43
1:A:263:ARG:NH2	1:A:278:TYR:HA	2.34	0.42
1:A:70:CYS:HB3	1:A:74:GLN:HE22	1.83	0.42
1:C:206:GLN:HE22	1:C:208:ASN:HB2	1.84	0.42
2:C:1001:ECL:CL8	2:C:1001:ECL:O20	2.74	0.42
1:D:422:GLY:O	1:D:467:TYR:OH	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:612:LEU:O	1:D:612:LEU:CG	2.67	0.42
1:B:496:LEU:HD21	1:D:475:LEU:HD11	2.01	0.42
2:D:1001:ECL:CL8	2:D:1001:ECL:O20	2.74	0.42
1:D:206:GLN:HE22	1:D:208:ASN:HB2	1.84	0.42
1:A:608:MET:HA	1:A:609:PRO:HD3	1.81	0.42
1:D:435:LEU:CD2	1:D:457:ALA:CB	2.81	0.42
1:D:70:CYS:HB3	1:D:74:GLN:HE22	1.83	0.42
2:A:1001:ECL:CL8	2:A:1001:ECL:O20	2.74	0.42
1:B:75:ARG:HH22	1:B:111:LEU:HD23	1.84	0.42
1:A:273:LEU:HD11	1:C:160:ILE:HG21	2.02	0.42
1:C:70:CYS:HB3	1:C:74:GLN:HE22	1.83	0.42
1:D:263:ARG:NH2	1:D:278:TYR:HA	2.34	0.42
1:B:263:ARG:NH2	1:B:278:TYR:HA	2.34	0.42
2:D:1001:ECL:CL8	2:D:1001:ECL:H8A	2.56	0.42
1:D:75:ARG:HH22	1:D:111:LEU:HD23	1.84	0.42
1:C:48:PRO:HA	1:C:51:GLN:HE22	1.84	0.42
1:A:612:LEU:CG	1:A:612:LEU:O	2.67	0.42
1:B:70:CYS:HB3	1:B:74:GLN:HE22	1.83	0.42
1:C:612:LEU:O	1:C:612:LEU:CG	2.68	0.42
1:A:48:PRO:HA	1:A:51:GLN:HE22	1.84	0.42
1:A:499:VAL:O	1:B:462:TRP:NE1	2.51	0.41
1:B:435:LEU:CD2	1:B:457:ALA:CB	2.81	0.41
1:C:554:MET:HE3	1:C:557:ILE:HD11	2.01	0.41
1:B:119:THR:HG22	1:B:122:HIS:CE1	2.56	0.41
1:B:237:ASN:OD1	1:B:240:GLY:N	2.53	0.41
1:B:422:GLY:O	1:B:467:TYR:OH	2.37	0.41
1:C:119:THR:HG22	1:C:122:HIS:CE1	2.56	0.41
1:D:160:ILE:HG21	1:C:273:LEU:HD11	2.03	0.41
1:B:435:LEU:HD12	1:B:435:LEU:C	2.38	0.41
1:D:119:THR:HG22	1:D:122:HIS:CE1	2.56	0.41
1:D:428:ILE:HG13	1:D:428:ILE:H	1.61	0.41
1:D:45:ARG:CG	1:D:45:ARG:NH1	2.72	0.41
1:A:204:ILE:HG21	1:A:247:ALA:HB2	2.03	0.41
1:D:202:ILE:HG13	1:D:202:ILE:H	1.71	0.41
1:B:121:LEU:HD12	1:B:122:HIS:N	2.36	0.41
1:D:496:LEU:HD21	1:C:475:LEU:HD11	2.03	0.41
1:D:121:LEU:HD12	1:D:122:HIS:N	2.36	0.41
1:B:114:PRO:HA	1:B:150:ALA:HB3	2.02	0.41
1:C:460:LEU:HD11	2:C:1001:ECL:C13	2.51	0.41
1:C:204:ILE:HG21	1:C:247:ALA:HB2	2.03	0.41
1:B:612:LEU:O	1:B:612:LEU:CG	2.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:460:LEU:HD11	2:D:1001:ECL:C13	2.51	0.41
1:B:460:LEU:HD11	2:B:1001:ECL:C13	2.51	0.40
1:D:160:ILE:HG13	1:D:160:ILE:H	1.79	0.40
1:A:237:ASN:OD1	1:A:241:LEU:N	2.47	0.40
1:C:422:GLY:O	1:C:467:TYR:OH	2.37	0.40
1:A:460:LEU:HD11	2:A:1001:ECL:C13	2.51	0.40
1:C:114:PRO:HA	1:C:150:ALA:HB3	2.02	0.40
1:A:119:THR:HG22	1:A:122:HIS:CE1	2.56	0.40
1:A:532:SER:O	1:A:536:LEU:HB2	2.22	0.40
1:C:121:LEU:HD12	1:C:122:HIS:N	2.36	0.40
1:A:121:LEU:HD12	1:A:122:HIS:N	2.36	0.40
1:A:114:PRO:HA	1:A:150:ALA:HB3	2.02	0.40
1:B:532:SER:O	1:B:536:LEU:HB2	2.22	0.40
1:D:114:PRO:HA	1:D:150:ALA:HB3	2.02	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 PDB entries followed by that with respect to all EM entries.

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	585/730 (80%)	551 (94%)	34 (6%)	0	100	100
1	B	585/730 (80%)	551 (94%)	34 (6%)	0	100	100
1	C	585/730 (80%)	551 (94%)	34 (6%)	0	100	100
1	D	585/730 (80%)	550 (94%)	35 (6%)	0	100	100
All	All	2340/2920 (80%)	2203 (94%)	137 (6%)	0	100	100

There are no Ramachandran outliers to report.

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 PDB entries followed by that with respect to all EM entries.

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	507/631 (80%)	488 (96%)	19 (4%)	39	68
1	B	507/631 (80%)	488 (96%)	19 (4%)	39	68
1	C	507/631 (80%)	488 (96%)	19 (4%)	39	68
1	D	507/631 (80%)	488 (96%)	19 (4%)	39	68
All	All	2028/2524 (80%)	1952 (96%)	76 (4%)	43	68

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	VAL
1	A	118	GLN
1	A	195	LEU
1	A	241	LEU
1	A	250	GLU
1	A	263	ARG
1	A	340	MET
1	A	400	LEU
1	A	413	SER
1	A	440	MET
1	A	441	VAL
1	A	442	MET
1	A	452	VAL
1	A	474	MET
1	A	492	ARG
1	A	497	MET
1	A	574	PHE
1	A	584	ARG
1	A	587	GLN
1	B	36	VAL
1	B	118	GLN
1	B	195	LEU
1	B	241	LEU
1	B	250	GLU

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Mol	Chain	Res	Type
1	B	263	ARG
1	B	340	MET
1	B	400	LEU
1	B	413	SER
1	B	440	MET
1	B	441	VAL
1	B	442	MET
1	B	452	VAL
1	B	474	MET
1	B	492	ARG
1	B	497	MET
1	B	574	PHE
1	B	584	ARG
1	B	587	GLN
1	D	36	VAL
1	D	118	GLN
1	D	195	LEU
1	D	241	LEU
1	D	250	GLU
1	D	263	ARG
1	D	340	MET
1	D	400	LEU
1	D	413	SER
1	D	440	MET
1	D	441	VAL
1	D	442	MET
1	D	452	VAL
1	D	474	MET
1	D	492	ARG
1	D	497	MET
1	D	574	PHE
1	D	584	ARG
1	D	587	GLN
1	C	36	VAL
1	C	118	GLN
1	C	195	LEU
1	C	241	LEU
1	C	250	GLU
1	C	263	ARG
1	C	340	MET
1	C	400	LEU
1	C	413	SER

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Mol	Chain	Res	Type
1	C	440	MET
1	C	441	VAL
1	C	442	MET
1	C	452	VAL
1	C	474	MET
1	C	492	ARG
1	C	497	MET
1	C	574	PHE
1	C	584	ARG
1	C	587	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	197	ASN
1	A	201	HIS
1	A	239	GLN
1	A	483	GLN
1	A	572	ASN
1	B	73	GLN
1	B	197	ASN
1	B	201	HIS
1	B	239	GLN
1	B	483	GLN
1	B	572	ASN
1	D	73	GLN
1	D	197	ASN
1	D	201	HIS
1	D	239	GLN
1	D	483	GLN
1	D	572	ASN
1	C	73	GLN
1	C	197	ASN
1	C	201	HIS
1	C	239	GLN
1	C	483	GLN
1	C	572	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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
2	ECL	A	1001	-	24,26,26	4.78	14 (58%)	32,35,35	0.96	2 (6%)
2	ECL	B	1001	-	24,26,26	4.79	14 (58%)	32,35,35	0.95	2 (6%)
2	ECL	C	1001	-	24,26,26	4.78	14 (58%)	32,35,35	0.96	2 (6%)
2	ECL	D	1001	-	24,26,26	4.79	14 (58%)	32,35,35	0.96	2 (6%)

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	ECL	A	1001	-	-	0/13/13/13	0/3/3/3
2	ECL	B	1001	-	-	0/13/13/13	0/3/3/3
2	ECL	C	1001	-	-	0/13/13/13	0/3/3/3
2	ECL	D	1001	-	-	0/13/13/13	0/3/3/3

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1001	ECL	C21-C17	-5.99	1.31	1.39
2	A	1001	ECL	C21-C17	-5.93	1.32	1.39
2	B	1001	ECL	C21-C17	-5.92	1.32	1.39
2	C	1001	ECL	C21-C17	-5.90	1.32	1.39
2	C	1001	ECL	C15-C16	-4.09	1.31	1.38
2	A	1001	ECL	C15-C16	-4.09	1.31	1.38
2	B	1001	ECL	C15-C16	-4.08	1.31	1.38
2	D	1001	ECL	C15-C16	-4.08	1.31	1.38
2	B	1001	ECL	C10-C9	-3.76	1.32	1.38
2	D	1001	ECL	C10-C9	-3.75	1.32	1.38
2	A	1001	ECL	C10-C9	-3.74	1.32	1.38
2	C	1001	ECL	C10-C9	-3.72	1.32	1.38
2	B	1001	ECL	C2-C1	-3.36	1.31	1.38
2	C	1001	ECL	C2-C1	-3.36	1.31	1.38
2	A	1001	ECL	C2-C1	-3.35	1.31	1.38
2	D	1001	ECL	C2-C1	-3.33	1.31	1.38
2	C	1001	ECL	C5-C14	-3.26	1.31	1.38
2	D	1001	ECL	C5-C14	-3.25	1.31	1.38
2	A	1001	ECL	C5-C14	-3.24	1.31	1.38
2	B	1001	ECL	C5-C14	-3.23	1.31	1.38
2	C	1001	ECL	C13-C11	-3.12	1.32	1.38
2	A	1001	ECL	C13-C11	-3.10	1.32	1.38
2	D	1001	ECL	C13-C11	-3.09	1.32	1.38
2	B	1001	ECL	C13-C11	-3.06	1.32	1.38
2	C	1001	ECL	C19-N1	-2.53	1.45	1.48
2	A	1001	ECL	C19-N1	-2.52	1.45	1.48
2	B	1001	ECL	C19-N1	-2.51	1.45	1.48
2	D	1001	ECL	C19-N1	-2.44	1.45	1.48
2	C	1001	ECL	C16-CL8	2.05	1.78	1.73
2	D	1001	ECL	C16-CL8	2.06	1.78	1.73
2	A	1001	ECL	C16-CL8	2.07	1.78	1.73
2	B	1001	ECL	C16-CL8	2.07	1.78	1.73
2	C	1001	ECL	C10-C1	6.79	1.52	1.38
2	A	1001	ECL	C10-C1	6.82	1.53	1.38
2	B	1001	ECL	C10-C1	6.83	1.53	1.38
2	D	1001	ECL	C10-C1	6.85	1.53	1.38
2	D	1001	ECL	C9-C11	7.71	1.53	1.38
2	A	1001	ECL	C9-C11	7.73	1.53	1.38
2	C	1001	ECL	C9-C11	7.76	1.53	1.38
2	B	1001	ECL	C9-C11	7.76	1.53	1.38
2	C	1001	ECL	C21-C5	7.84	1.52	1.38
2	A	1001	ECL	C21-C5	7.86	1.52	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1001	ECL	C21-C5	7.87	1.52	1.38
2	B	1001	ECL	C21-C5	7.87	1.52	1.38
2	C	1001	ECL	C13-C2	7.97	1.53	1.38
2	B	1001	ECL	C13-C2	7.97	1.53	1.38
2	A	1001	ECL	C13-C2	8.01	1.53	1.38
2	D	1001	ECL	C13-C2	8.03	1.53	1.38
2	C	1001	ECL	C15-C14	8.47	1.52	1.38
2	A	1001	ECL	C15-C14	8.50	1.52	1.38
2	B	1001	ECL	C15-C14	8.51	1.52	1.38
2	D	1001	ECL	C15-C14	8.53	1.52	1.38
2	A	1001	ECL	C16-C17	11.05	1.53	1.39
2	C	1001	ECL	C16-C17	11.06	1.53	1.39
2	D	1001	ECL	C16-C17	11.07	1.53	1.39
2	B	1001	ECL	C16-C17	11.10	1.53	1.39

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1001	ECL	C15-C16-C17	-2.06	119.99	122.44
2	D	1001	ECL	C15-C16-C17	-2.05	120.00	122.44
2	B	1001	ECL	C15-C16-C17	-2.04	120.01	122.44
2	A	1001	ECL	C15-C16-C17	-2.03	120.02	122.44
2	B	1001	ECL	C21-C17-C16	2.37	119.24	116.80
2	A	1001	ECL	C21-C17-C16	2.38	119.24	116.80
2	D	1001	ECL	C21-C17-C16	2.38	119.25	116.80
2	C	1001	ECL	C21-C17-C16	2.38	119.25	116.80

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 32 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	ECL	8	0
2	B	1001	ECL	8	0
2	C	1001	ECL	8	0
2	D	1001	ECL	8	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.