



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:21 PM GMT

PDB ID : 4I1P
Title : Human MALT1 (caspase-IG3) in complex with activity based-probe
Authors : Schlauderer, F.; Lammens, K.; Hopfner, K.P.
Deposited on : 2012-11-21
Resolution : 2.40 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

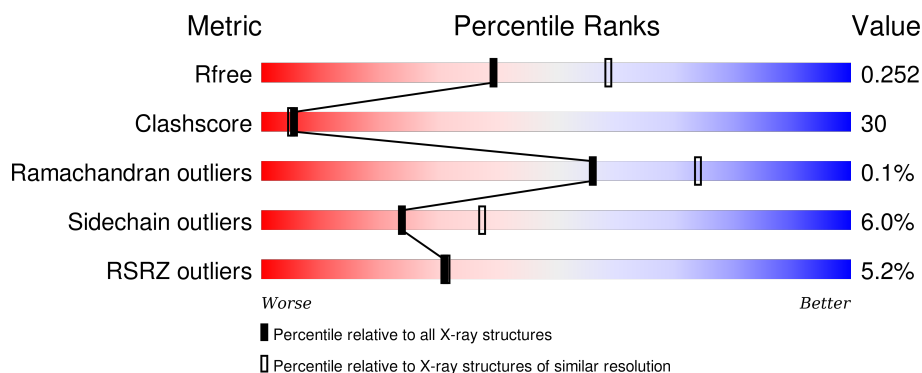
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	388	<div> <div>5%</div> <div>63%</div> <div>30%</div> <div>• •</div> </div>
1	C	388	<div> <div>5%</div> <div>62%</div> <div>32%</div> <div>• •</div> </div>
2	B	5	<div> <div>40%</div> <div>60%</div> </div>
2	D	5	<div> <div>40%</div> <div>60%</div> </div>

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
3	ACT	A	801	-	-	X	X
3	ACT	A	803	-	-	X	X
3	ACT	A	804	-	-	X	X
3	ACT	A	805	-	-	-	X
3	ACT	C	801	-	-	-	X
3	ACT	C	803	-	-	-	X
3	ACT	C	805	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6256 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 Mucosa-associated lymphoid tissue lymphoma translocation protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	124	0	0
			3007	1925	487	574	21			
1	C	380	Total	C	N	O	S	113	0	0
			3015	1930	488	575	22			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	338	MET	-	INITIATING METHIONINE	UNP Q9UDY8
A	720	HIS	-	EXPRESSION TAG	UNP Q9UDY8
A	721	HIS	-	EXPRESSION TAG	UNP Q9UDY8
A	722	HIS	-	EXPRESSION TAG	UNP Q9UDY8
A	723	HIS	-	EXPRESSION TAG	UNP Q9UDY8
A	724	HIS	-	EXPRESSION TAG	UNP Q9UDY8
A	725	HIS	-	EXPRESSION TAG	UNP Q9UDY8
C	338	MET	-	EXPRESSION TAG	UNP Q9UDY8
C	720	HIS	-	EXPRESSION TAG	UNP Q9UDY8
C	721	HIS	-	EXPRESSION TAG	UNP Q9UDY8
C	722	HIS	-	EXPRESSION TAG	UNP Q9UDY8
C	723	HIS	-	EXPRESSION TAG	UNP Q9UDY8
C	724	HIS	-	EXPRESSION TAG	UNP Q9UDY8
C	725	HIS	-	EXPRESSION TAG	UNP Q9UDY8

- Molecule 2 is a protein called tetrapeptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	5	Total	C	N	O	0	0	1
			39	23	10	6			
2	D	5	Total	C	N	O	0	0	1
			39	23	10	6			

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	58	Total	O	0	0
			58	58		

Continued on next page...

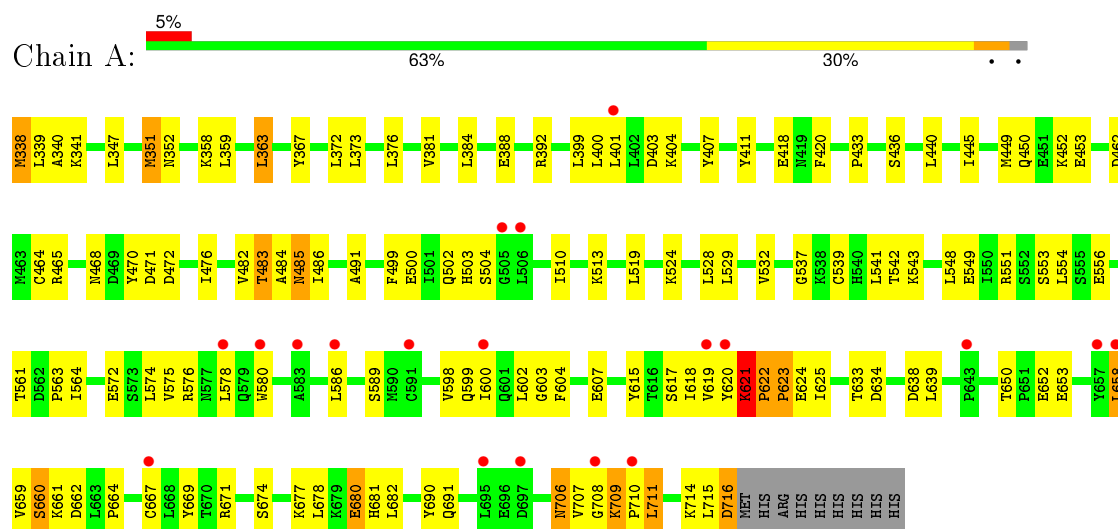
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total 2	O 2	0	0
4	C	55	Total 55	O 55	0	0
4	D	1	Total 1	O 1	0	0

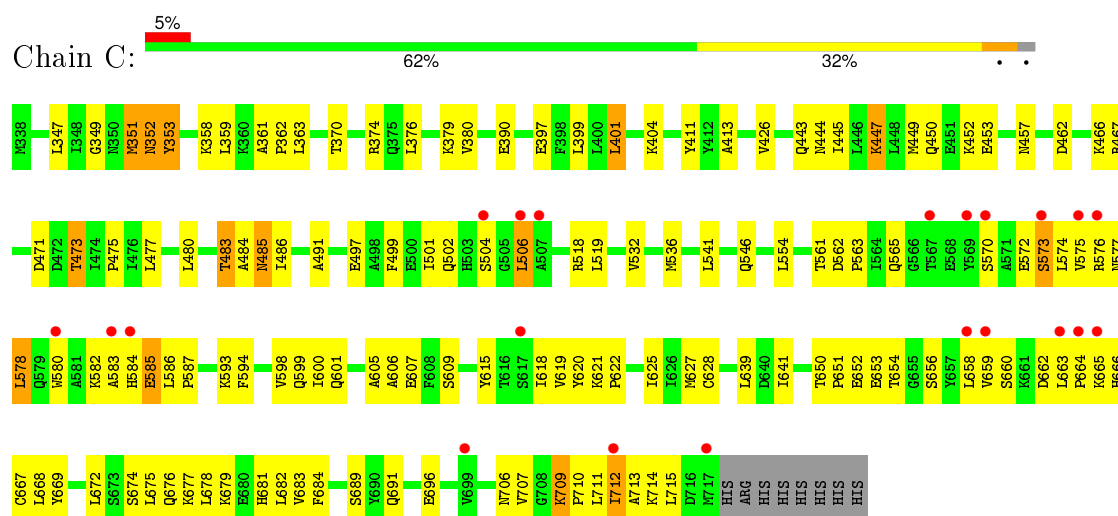
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Mucosa-associated lymphoid tissue lymphoma translocation protein 1



- Molecule 1: Mucosa-associated lymphoid tissue lymphoma translocation protein 1



- Molecule 2: tetrapeptide



A0	L1	R2	S3	?4
----	----	----	----	----

- Molecule 2: tetrapeptide

Chain D:



A0	L1	R2	S3	?4
----	----	----	----	----

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.24Å 53.48Å 78.74Å 90.00° 103.80° 90.00°	Depositor
Resolution (Å)	43.94 – 2.40 43.94 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.1 (43.94-2.40) 96.3 (43.94-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.198 , 0.245 0.205 , 0.252	Depositor DCC
R_{free} test set	1370 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	52.2	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 57.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 27391 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6256	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 4AR, ACT

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	0.57	2/3062 (0.1%)	0.64	1/4141 (0.0%)
1	C	0.52	0/3070	0.61	1/4151 (0.0%)
2	B	0.18	0/26	0.25	0/32
2	D	0.25	0/26	0.48	0/32
All	All	0.54	2/6184 (0.0%)	0.63	2/8356 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	623	PRO	N-CD	5.21	1.55	1.47
1	A	622	PRO	N-CD	5.16	1.55	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	621	LYS	C-N-CD	5.74	140.46	128.40
1	C	352	ASN	C-N-CA	-5.58	107.75	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3007	0	3029	152	0
1	C	3015	0	3037	188	0
2	B	39	0	41	8	0
2	D	39	0	41	19	0
3	A	20	0	15	12	0
3	C	20	0	15	2	0
4	A	58	0	0	8	0
4	B	2	0	0	0	0
4	C	55	0	0	3	0
4	D	1	0	0	0	0
All	All	6256	0	6178	347	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 30.

All (347) 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)
2:D:4:4AR:C2	2:D:4:4AR:C7	1.92	1.45
1:A:483:THR:HB	1:A:485:ASN:ND2	1.40	1.35
1:C:473:THR:CG2	1:C:475:PRO:HD3	1.61	1.29
1:A:464:CYS:CB	2:B:4:4AR:C7	2.17	1.22
1:C:572:GLU:O	1:C:576:ARG:HG3	1.38	1.21
1:A:551:ARG:CD	3:A:801:ACT:H1	1.73	1.18
1:C:363:LEU:CD2	1:C:658:LEU:HD11	1.75	1.14
1:C:607:GLU:HG3	1:C:659:VAL:HG23	1.18	1.09
1:C:363:LEU:CD2	1:C:658:LEU:CD1	2.31	1.09
1:A:483:THR:CB	1:A:485:ASN:HD21	1.66	1.08
1:C:473:THR:HG23	1:C:475:PRO:HD3	1.18	1.07
1:A:551:ARG:HD3	3:A:801:ACT:CH3	1.84	1.07
1:C:578:LEU:CD1	1:C:582:LYS:NZ	2.20	1.05
1:C:363:LEU:HD23	1:C:658:LEU:HD11	1.08	1.05
1:C:627:MET:HE1	1:C:691:GLN:NE2	1.72	1.05
1:A:401:LEU:HD21	1:A:580:TRP:CE3	1.91	1.03
1:A:551:ARG:HD3	3:A:801:ACT:H1	1.06	1.03
2:D:4:4AR:C1	2:D:4:4AR:C7	2.37	1.02
1:A:401:LEU:CD2	1:A:580:TRP:CE3	2.44	1.00
1:C:541:LEU:HG	2:D:1:LEU:HD11	1.42	1.00
1:C:363:LEU:HD23	1:C:658:LEU:CD1	1.92	0.99
1:A:551:ARG:CD	3:A:801:ACT:CH3	2.39	0.98
1:A:483:THR:CB	1:A:485:ASN:ND2	2.24	0.97
1:C:587:PRO:HB3	1:C:707:VAL:CG1	1.93	0.97

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:587:PRO:CG	1:C:707:VAL:HG11	1.95	0.97
4:C:930:HOH:O	2:D:2:ARG:HD2	1.65	0.96
1:C:363:LEU:HD21	1:C:658:LEU:CD1	1.93	0.95
1:A:483:THR:HB	1:A:485:ASN:HD21	1.14	0.95
1:A:401:LEU:HD21	1:A:580:TRP:HE3	1.25	0.94
1:C:483:THR:HB	1:C:485:ASN:ND2	1.83	0.94
1:C:484:ALA:O	1:C:554:LEU:HA	1.69	0.93
1:C:352:ASN:HB2	1:C:358:LYS:HE3	1.50	0.93
1:C:351:MET:HG2	1:C:359:LEU:O	1.70	0.92
1:C:578:LEU:HD12	1:C:582:LYS:NZ	1.83	0.92
1:A:681:HIS:HB3	1:A:708:GLY:HA2	1.48	0.92
1:C:615:TYR:CE2	1:C:664:PRO:HG2	2.05	0.91
1:C:676:GLN:HB2	3:C:803:ACT:H2	1.53	0.89
1:C:587:PRO:HB3	1:C:707:VAL:HG13	1.52	0.89
1:C:374:ARG:CZ	1:C:584:HIS:HE1	1.84	0.89
1:C:615:TYR:HE2	1:C:664:PRO:HG2	1.35	0.88
1:C:353:TYR:HE2	1:C:426:VAL:O	1.55	0.88
1:A:572:GLU:O	1:A:575:VAL:HG12	1.74	0.88
1:C:541:LEU:CG	2:D:1:LEU:HD11	2.03	0.88
1:C:587:PRO:HG3	1:C:707:VAL:HG11	1.53	0.87
1:C:473:THR:HG22	1:C:475:PRO:HD3	1.54	0.87
1:C:586:LEU:HD11	1:C:606:ALA:N	1.91	0.86
1:C:578:LEU:HD12	1:C:582:LYS:HZ3	1.41	0.86
1:C:578:LEU:CD1	1:C:582:LYS:HZ2	1.84	0.86
1:C:587:PRO:CB	1:C:707:VAL:HG11	2.04	0.86
1:C:578:LEU:CD1	1:C:582:LYS:HZ3	1.86	0.85
2:D:4:4AR:N1	2:D:4:4AR:C7	2.40	0.85
1:A:586:LEU:HD13	1:A:604:PHE:HB2	1.57	0.85
1:A:352:ASN:HB2	4:A:915:HOH:O	1.76	0.85
1:A:464:CYS:HG	2:B:4:4AR:C7	0.89	0.84
1:C:627:MET:CE	1:C:691:GLN:NE2	2.40	0.84
1:C:583:ALA:O	1:C:711:LEU:HA	1.76	0.84
1:A:667:CYS:O	1:A:669:TYR:CE2	2.30	0.84
1:C:587:PRO:CB	1:C:707:VAL:CG1	2.56	0.83
1:C:353:TYR:CE2	1:C:426:VAL:O	2.30	0.83
1:A:603:GLY:HA3	1:A:615:TYR:CE1	2.14	0.83
1:C:627:MET:HE2	1:C:691:GLN:HG3	1.60	0.83
1:A:352:ASN:OD1	1:A:358:LYS:HD3	1.79	0.82
1:C:572:GLU:O	1:C:576:ARG:CG	2.26	0.82
1:C:541:LEU:CD2	2:D:1:LEU:HD11	2.10	0.81
1:A:468:ASN:HB2	4:A:957:HOH:O	1.81	0.81

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:401:LEU:CD2	1:A:580:TRP:CZ3	2.63	0.80
1:A:660:SER:C	1:A:662:ASP:H	1.83	0.80
1:A:381:VAL:HG23	1:A:580:TRP:CZ2	2.16	0.80
1:C:707:VAL:HG12	1:C:707:VAL:O	1.80	0.80
1:A:367:TYR:CB	1:A:658:LEU:HD21	2.12	0.78
1:A:483:THR:HB	1:A:485:ASN:CG	2.04	0.77
1:C:607:GLU:HG3	1:C:659:VAL:CG2	2.08	0.77
1:A:401:LEU:HD23	1:A:580:TRP:CZ3	2.19	0.76
1:C:586:LEU:HD23	1:C:711:LEU:CD1	2.15	0.76
1:C:678:LEU:O	1:C:678:LEU:HD12	1.84	0.76
1:C:483:THR:HB	1:C:485:ASN:CG	2.05	0.76
1:A:598:VAL:HG13	1:A:620:TYR:O	1.86	0.75
2:D:4:4AR:O1	2:D:4:4AR:C7	2.35	0.75
1:C:374:ARG:CZ	1:C:584:HIS:CE1	2.70	0.74
1:C:473:THR:HG23	1:C:475:PRO:CD	2.10	0.74
1:A:551:ARG:HD2	3:A:801:ACT:CH3	2.17	0.74
1:C:627:MET:HE1	1:C:691:GLN:CD	2.07	0.73
1:A:401:LEU:HG	1:A:580:TRP:CZ3	2.23	0.73
1:A:450:GLN:NE2	1:A:483:THR:OG1	2.20	0.73
1:A:681:HIS:CD2	1:A:708:GLY:HA3	2.23	0.73
1:C:447:LYS:HE2	1:C:480:LEU:HD23	1.71	0.73
1:A:681:HIS:CB	1:A:708:GLY:HA2	2.20	0.72
1:A:681:HIS:CD2	1:A:708:GLY:CA	2.72	0.72
1:A:483:THR:CG2	1:A:485:ASN:OD1	2.37	0.71
1:A:420:PHE:CE2	4:A:949:HOH:O	2.43	0.71
1:C:363:LEU:HD21	1:C:658:LEU:HD12	1.72	0.71
1:C:627:MET:CE	1:C:691:GLN:CD	2.59	0.71
1:C:578:LEU:HD11	1:C:582:LYS:NZ	2.06	0.71
1:C:609:SER:O	1:C:712:ILE:HG22	1.91	0.70
1:C:483:THR:HB	1:C:485:ASN:HD21	1.55	0.70
1:C:473:THR:CG2	1:C:475:PRO:CD	2.56	0.70
1:C:627:MET:HE2	1:C:691:GLN:CG	2.22	0.70
1:C:483:THR:CB	1:C:485:ASN:HD21	2.04	0.69
1:A:660:SER:C	1:A:662:ASP:N	2.43	0.69
1:A:617:SER:HB3	1:A:667:CYS:SG	2.33	0.68
1:C:397:GLU:HG2	1:C:715:LEU:HD13	1.74	0.68
1:C:593:LYS:HG3	1:C:599:GLN:HG2	1.73	0.68
1:A:404:LYS:HA	1:A:453:GLU:O	1.93	0.68
1:A:551:ARG:NH2	4:A:923:HOH:O	2.23	0.67
1:C:600:ILE:HG22	1:C:618:ILE:HA	1.76	0.67
1:C:615:TYR:CE2	1:C:664:PRO:CG	2.77	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:501:ILE:HG22	1:C:504:SER:HB3	1.77	0.67
1:C:674:SER:HB3	1:C:677:LYS:HG3	1.75	0.67
1:C:401:LEU:HD13	1:C:580:TRP:CE3	2.30	0.66
1:C:682:LEU:HD23	1:C:684:PHE:HD1	1.60	0.66
1:A:551:ARG:HD2	3:A:801:ACT:H1	1.73	0.66
1:A:711:LEU:O	1:A:714:LYS:HG3	1.95	0.66
1:C:712:ILE:O	1:C:712:ILE:HD13	1.96	0.66
1:C:586:LEU:HD23	1:C:711:LEU:HD13	1.76	0.65
1:C:483:THR:CB	1:C:485:ASN:ND2	2.58	0.65
1:A:621:LYS:HD3	1:A:625:ILE:O	1.96	0.65
1:A:420:PHE:HE2	4:A:949:HOH:O	1.79	0.65
1:C:675:LEU:O	1:C:678:LEU:HG	1.97	0.65
1:C:586:LEU:HD23	1:C:711:LEU:HD11	1.78	0.64
1:C:363:LEU:HD21	1:C:658:LEU:CG	2.26	0.64
1:C:659:VAL:O	1:C:659:VAL:HG12	1.95	0.64
1:C:615:TYR:CE2	1:C:664:PRO:HD2	2.32	0.64
1:C:682:LEU:CD2	1:C:684:PHE:CD1	2.81	0.64
1:A:401:LEU:HD23	1:A:580:TRP:HZ3	1.63	0.64
1:A:681:HIS:HD2	1:A:708:GLY:CA	2.10	0.64
1:A:338:MET:HA	1:A:338:MET:HE3	1.80	0.64
1:C:598:VAL:HG13	1:C:620:TYR:O	1.98	0.64
1:C:374:ARG:NH2	1:C:584:HIS:HE1	1.95	0.64
1:A:681:HIS:HD2	1:A:708:GLY:HA3	1.63	0.63
1:A:551:ARG:CD	3:A:801:ACT:H2	2.29	0.63
1:C:666:HIS:O	1:C:667:CYS:SG	2.54	0.63
1:C:682:LEU:HD23	1:C:684:PHE:CD1	2.34	0.63
1:A:706:ASN:O	1:A:706:ASN:ND2	2.32	0.63
1:A:339:LEU:CD2	1:A:563:PRO:HB2	2.29	0.63
1:C:712:ILE:C	1:C:712:ILE:HD13	2.19	0.62
1:C:352:ASN:HB3	4:C:907:HOH:O	1.99	0.62
1:C:605:ALA:CB	1:C:663:LEU:HD11	2.30	0.62
1:A:351:MET:HB3	1:A:359:LEU:O	1.99	0.62
1:A:483:THR:HG22	1:A:485:ASN:OD1	1.98	0.62
1:C:363:LEU:CD2	1:C:658:LEU:CG	2.78	0.62
1:C:628:CYS:HA	1:C:689:SER:O	1.99	0.62
1:C:361:ALA:N	1:C:362:PRO:CD	2.63	0.61
1:C:594:PHE:CE1	1:C:600:ILE:HD11	2.34	0.61
1:A:715:LEU:O	1:A:716:ASP:CB	2.47	0.61
1:C:363:LEU:HD21	1:C:658:LEU:HG	1.81	0.61
1:C:586:LEU:HD11	1:C:606:ALA:HB2	1.83	0.61
1:A:392:ARG:NH1	3:A:803:ACT:OXT	2.29	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:600:ILE:CD1	1:A:618:ILE:HD13	2.31	0.61
1:C:615:TYR:CE2	1:C:664:PRO:CD	2.84	0.61
1:A:572:GLU:O	1:A:575:VAL:CG1	2.49	0.61
1:A:401:LEU:CG	1:A:580:TRP:CZ3	2.84	0.60
1:C:379:LYS:HE2	1:C:577:ASN:ND2	2.16	0.60
1:C:709:LYS:O	1:C:714:LYS:HE3	2.01	0.60
1:C:501:ILE:CG2	1:C:504:SER:HB3	2.31	0.60
1:A:483:THR:O	1:A:484:ALA:HB3	2.00	0.60
1:C:541:LEU:HG	2:D:1:LEU:CD1	2.25	0.60
1:C:586:LEU:HD11	1:C:606:ALA:CB	2.31	0.60
1:A:682:LEU:HD23	1:A:707:VAL:HG22	1.83	0.60
1:C:586:LEU:CD2	1:C:711:LEU:HD11	2.32	0.60
1:A:621:LYS:CD	1:A:625:ILE:O	2.50	0.60
1:A:715:LEU:O	1:A:716:ASP:HB3	2.01	0.59
1:C:466:LYS:HE3	1:C:497:GLU:HG3	1.84	0.59
1:A:532:VAL:HG12	1:A:548:LEU:HD12	1.84	0.59
1:A:418:GLU:OE1	1:A:465:ARG:HD2	2.01	0.59
1:C:659:VAL:HG12	1:C:663:LEU:CD1	2.33	0.59
1:A:664:PRO:O	1:A:669:TYR:OH	2.21	0.59
1:C:641:ILE:HD11	1:C:672:LEU:HD22	1.85	0.59
1:C:707:VAL:CG1	1:C:707:VAL:O	2.50	0.59
1:C:601:GLN:HB2	1:C:619:VAL:CG2	2.34	0.58
1:C:483:THR:O	1:C:484:ALA:HB3	2.03	0.58
1:C:639:LEU:HD21	1:C:678:LEU:HD23	1.86	0.58
1:C:450:GLN:NE2	1:C:483:THR:OG1	2.34	0.58
1:A:603:GLY:HA3	1:A:615:TYR:CZ	2.38	0.58
1:C:715:LEU:O	3:C:805:ACT:H1	2.04	0.58
1:A:347:LEU:HG	1:A:411:TYR:HB3	1.86	0.57
1:A:464:CYS:HB2	2:B:4:4AR:C7	2.29	0.57
1:A:682:LEU:HD23	1:A:707:VAL:CG2	2.35	0.57
1:A:681:HIS:HD2	1:A:709:LYS:N	2.03	0.56
1:C:536:MET:CE	1:C:546:GLN:HG2	2.36	0.56
1:A:503:HIS:CE1	1:A:504:SER:O	2.58	0.56
1:C:682:LEU:CD2	1:C:684:PHE:CE1	2.89	0.56
1:C:607:GLU:CG	1:C:659:VAL:HG23	2.13	0.56
1:C:374:ARG:NH2	1:C:584:HIS:CE1	2.73	0.56
1:A:367:TYR:HB3	1:A:658:LEU:HD21	1.88	0.55
1:A:600:ILE:HD12	1:A:618:ILE:HD13	1.87	0.55
1:A:367:TYR:CG	1:A:658:LEU:CD2	2.89	0.55
1:C:578:LEU:HD11	1:C:582:LYS:HZ3	1.67	0.55
1:A:599:GLN:O	1:A:619:VAL:HG12	2.05	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:363:LEU:CD2	1:C:658:LEU:HG	2.37	0.55
1:A:543:LYS:HE3	4:A:932:HOH:O	2.07	0.54
1:A:484:ALA:O	1:A:554:LEU:HA	2.07	0.54
1:C:541:LEU:HD21	2:D:1:LEU:HD11	1.89	0.54
1:C:621:LYS:HE3	1:C:625:ILE:O	2.08	0.54
1:C:483:THR:HB	1:C:485:ASN:OD1	2.07	0.54
1:C:499:PHE:CG	2:D:1:LEU:HD23	2.43	0.54
1:C:351:MET:CG	1:C:359:LEU:O	2.50	0.54
1:A:551:ARG:HD2	3:A:801:ACT:H2	1.90	0.53
1:C:473:THR:HG22	1:C:475:PRO:CD	2.32	0.53
1:A:367:TYR:HB2	1:A:658:LEU:HD21	1.88	0.53
1:A:476:ILE:HG22	1:A:476:ILE:O	2.09	0.53
1:C:656:SER:O	1:C:660:SER:OG	2.26	0.53
1:C:347:LEU:HG	1:C:411:TYR:HB3	1.90	0.53
1:C:599:GLN:NE2	4:C:917:HOH:O	2.41	0.53
1:A:485:ASN:H	1:A:485:ASN:ND2	2.06	0.53
1:C:659:VAL:HG12	1:C:663:LEU:HD13	1.91	0.53
1:A:482:VAL:O	1:A:482:VAL:HG12	2.08	0.52
1:A:681:HIS:HD2	1:A:709:LYS:H	1.58	0.52
1:A:537:GLY:O	1:A:543:LYS:NZ	2.31	0.52
1:C:570:SER:OG	1:C:573:SER:HB2	2.09	0.52
1:A:660:SER:O	1:A:662:ASP:N	2.42	0.52
1:C:682:LEU:HD13	1:C:710:PRO:HG3	1.90	0.52
1:A:500:GLU:CG	2:B:2:ARG:HB3	2.40	0.52
1:A:639:LEU:HD21	1:A:678:LEU:HG	1.91	0.52
1:C:573:SER:HA	1:C:576:ARG:HD2	1.92	0.52
1:C:578:LEU:HD12	1:C:582:LYS:CE	2.39	0.52
1:A:367:TYR:CB	1:A:658:LEU:CD2	2.86	0.52
1:C:401:LEU:HD22	1:C:580:TRP:CD2	2.45	0.52
1:A:667:CYS:O	1:A:669:TYR:CD2	2.62	0.52
1:A:589:SER:HB2	1:A:602:LEU:O	2.09	0.52
1:C:578:LEU:HD13	1:C:582:LYS:HZ2	1.72	0.52
2:D:4:4AR:C7	2:D:4:4AR:C3	2.88	0.51
1:C:627:MET:HE2	1:C:691:GLN:CD	2.29	0.51
1:A:445:ILE:O	1:A:449:MET:HG3	2.11	0.51
1:C:363:LEU:CD2	1:C:658:LEU:HD12	2.34	0.51
1:C:390:GLU:OE2	1:C:677:LYS:HE3	2.11	0.51
1:C:518:ARG:HG3	1:C:532:VAL:HG22	1.92	0.51
1:A:659:VAL:HG13	1:A:662:ASP:HB3	1.93	0.50
1:A:618:ILE:HG22	1:A:618:ILE:O	2.11	0.50
1:C:445:ILE:O	1:C:449:MET:HG3	2.11	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:388:GLU:O	1:A:392:ARG:HG3	2.12	0.50
1:C:353:TYR:CD1	1:C:353:TYR:N	2.79	0.50
1:A:401:LEU:CG	1:A:580:TRP:CE3	2.95	0.50
1:C:585:GLU:O	1:C:711:LEU:HD12	2.11	0.50
1:A:399:LEU:O	1:A:452:LYS:NZ	2.39	0.50
1:A:674:SER:HB3	1:A:677:LYS:HD2	1.94	0.50
1:C:541:LEU:HD21	2:D:1:LEU:CD1	2.42	0.50
1:C:652:GLU:OE2	1:C:660:SER:HB3	2.11	0.50
1:A:500:GLU:HG3	2:B:2:ARG:HB3	1.94	0.50
1:C:678:LEU:HD12	1:C:678:LEU:C	2.31	0.49
1:C:682:LEU:HD21	1:C:684:PHE:CD1	2.47	0.49
1:C:502:GLN:HG3	2:D:2:ARG:NH1	2.27	0.49
1:A:625:ILE:HG21	1:A:690:TYR:HB2	1.94	0.49
1:A:485:ASN:HD22	1:A:486:ILE:HG13	1.78	0.49
1:C:352:ASN:HB3	1:C:358:LYS:HD3	1.95	0.49
1:C:609:SER:HB3	1:C:712:ILE:HB	1.95	0.49
1:A:662:ASP:O	1:A:662:ASP:OD1	2.30	0.49
1:A:499:PHE:HA	2:B:2:ARG:O	2.13	0.48
1:C:374:ARG:NE	1:C:584:HIS:CE1	2.81	0.48
1:C:471:ASP:OD1	1:C:473:THR:HB	2.14	0.48
1:A:524:LYS:HE3	1:A:556:GLU:O	2.14	0.48
1:A:680:GLU:O	1:A:681:HIS:C	2.51	0.48
1:A:625:ILE:HG23	1:A:691:GLN:O	2.14	0.48
1:A:376:LEU:HD22	1:A:561:THR:HG22	1.96	0.48
1:A:367:TYR:CG	1:A:658:LEU:HD21	2.48	0.48
1:A:339:LEU:HD22	1:A:563:PRO:HB2	1.95	0.48
1:A:650:THR:OG1	1:A:653:GLU:HG3	2.13	0.48
1:C:586:LEU:CD1	1:C:606:ALA:N	2.72	0.47
1:C:361:ALA:HB3	2:D:4:4AR:HH11	1.79	0.47
1:A:367:TYR:HB2	1:A:658:LEU:CD2	2.45	0.47
1:C:578:LEU:O	1:C:582:LYS:HD3	2.15	0.47
1:C:659:VAL:CG1	1:C:659:VAL:O	2.61	0.47
1:A:392:ARG:HB3	3:A:803:ACT:H1	1.96	0.47
1:A:338:MET:HA	1:A:338:MET:CE	2.43	0.47
1:C:370:THR:HG23	1:C:380:VAL:HG11	1.97	0.47
1:C:502:GLN:HG3	2:D:2:ARG:HH12	1.80	0.46
1:A:392:ARG:HH12	3:A:804:ACT:C	2.28	0.46
1:A:510:ILE:O	1:A:513:LYS:HB3	2.14	0.46
1:A:622:PRO:HB2	1:A:624:GLU:CD	2.35	0.46
1:C:574:LEU:O	1:C:578:LEU:HB2	2.16	0.46
1:A:574:LEU:O	1:A:578:LEU:HD23	2.16	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:513:LYS:HG2	4:A:950:HOH:O	2.14	0.46
1:C:651:PRO:O	1:C:654:THR:OG1	2.30	0.46
1:C:379:LYS:HE2	1:C:577:ASN:HD21	1.80	0.45
1:C:681:HIS:CE1	1:C:709:LYS:HD2	2.52	0.45
1:A:529:LEU:O	1:A:532:VAL:HB	2.16	0.45
1:C:404:LYS:HA	1:C:453:GLU:O	2.17	0.45
1:C:609:SER:O	1:C:712:ILE:CG2	2.61	0.45
1:A:660:SER:O	1:A:661:LYS:HB2	2.16	0.45
1:C:572:GLU:HA	1:C:575:VAL:HG12	1.99	0.45
1:C:587:PRO:CB	1:C:707:VAL:HG13	2.29	0.45
1:C:359:LEU:HD23	2:D:2:ARG:HD3	1.99	0.45
1:A:607:GLU:OE2	1:A:671:ARG:NH2	2.46	0.45
1:A:681:HIS:CG	1:A:708:GLY:HA2	2.52	0.44
1:A:575:VAL:HG13	1:A:576:ARG:N	2.31	0.44
1:A:433:PRO:HG2	1:A:470:TYR:CE1	2.52	0.44
1:C:650:THR:OG1	1:C:653:GLU:HG3	2.17	0.44
1:A:483:THR:O	1:A:484:ALA:CB	2.66	0.44
1:C:572:GLU:O	1:C:575:VAL:HG13	2.16	0.44
1:A:381:VAL:HG23	1:A:580:TRP:CE2	2.52	0.44
1:C:506:LEU:H	1:C:506:LEU:HG	1.54	0.44
1:A:401:LEU:HG	1:A:580:TRP:CE3	2.52	0.44
1:C:601:GLN:HB2	1:C:619:VAL:HG23	1.98	0.44
1:C:353:TYR:N	1:C:353:TYR:HD1	2.15	0.44
1:A:528:LEU:O	1:A:532:VAL:HG23	2.18	0.44
2:D:4:4AR:HB3	2:D:4:4AR:C7	2.48	0.44
1:C:594:PHE:HE1	1:C:600:ILE:HD11	1.77	0.44
1:C:586:LEU:HD11	1:C:606:ALA:CA	2.48	0.44
1:C:462:ASP:HA	1:C:491:ALA:HB2	2.00	0.44
1:A:652:GLU:N	1:A:652:GLU:OE1	2.42	0.43
1:C:682:LEU:HD21	1:C:684:PHE:CE1	2.53	0.43
1:C:598:VAL:HG22	1:C:622:PRO:HD2	2.00	0.43
1:C:519:LEU:HG	1:C:519:LEU:O	2.19	0.43
1:C:352:ASN:CB	1:C:358:LYS:HE3	2.35	0.43
1:C:443:GLN:HB3	1:C:477:LEU:HD22	2.00	0.43
1:A:539:CYS:HB3	1:A:542:THR:OG1	2.18	0.43
1:C:683:VAL:HG22	1:C:706:ASN:OD1	2.19	0.43
1:C:583:ALA:HB1	1:C:712:ILE:HA	2.01	0.43
1:C:536:MET:HE1	1:C:546:GLN:HG2	2.01	0.43
1:A:372:LEU:HB3	1:A:519:LEU:HD23	2.01	0.43
1:A:363:LEU:HD12	1:A:384:LEU:HD11	2.00	0.42
1:A:340:ALA:HB2	1:A:407:TYR:CE1	2.54	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:483:THR:O	1:C:484:ALA:CB	2.68	0.42
1:C:586:LEU:CD2	1:C:711:LEU:CD1	2.90	0.42
1:A:340:ALA:HB2	1:A:407:TYR:CZ	2.54	0.42
1:A:440:LEU:HA	3:A:804:ACT:O	2.20	0.42
1:C:586:LEU:HD11	1:C:606:ALA:H	1.77	0.42
1:A:633:THR:OG1	1:A:634:ASP:N	2.52	0.42
1:C:376:LEU:HD22	1:C:561:THR:HG22	2.01	0.42
1:C:562:ASP:HA	1:C:563:PRO:HD3	1.83	0.42
1:A:623:PRO:O	1:A:624:GLU:C	2.58	0.42
1:A:401:LEU:CD2	1:A:580:TRP:HE3	1.98	0.42
1:C:499:PHE:CB	2:D:1:LEU:HD23	2.49	0.42
1:A:599:GLN:O	1:A:619:VAL:CG1	2.68	0.42
1:A:400:LEU:HA	1:A:400:LEU:HD23	1.81	0.42
1:C:457:ASN:ND2	1:C:485:ASN:ND2	2.67	0.42
1:A:706:ASN:ND2	1:A:706:ASN:C	2.73	0.42
1:A:715:LEU:O	1:A:716:ASP:OD1	2.37	0.41
1:C:651:PRO:HD3	1:C:669:TYR:CE2	2.55	0.41
1:C:583:ALA:HA	1:C:711:LEU:O	2.20	0.41
1:A:681:HIS:CD2	1:A:708:GLY:HA2	2.55	0.41
1:A:658:LEU:HA	1:A:658:LEU:HD12	1.81	0.41
1:C:349:GLY:HA2	1:C:413:ALA:O	2.19	0.41
1:A:621:LYS:HD2	1:A:625:ILE:O	2.19	0.41
1:C:615:TYR:CZ	1:C:664:PRO:CD	3.03	0.41
1:C:379:LYS:HB3	1:C:580:TRP:CD1	2.56	0.41
1:A:602:LEU:HD12	1:A:602:LEU:HA	1.74	0.41
1:A:436:SER:HB3	1:A:471:ASP:OD2	2.21	0.41
1:A:373:LEU:HA	1:A:373:LEU:HD23	1.86	0.41
1:C:681:HIS:HE1	1:C:709:LYS:HD2	1.86	0.41
1:C:712:ILE:HG23	1:C:713:ALA:N	2.36	0.41
1:A:500:GLU:OE2	2:B:2:ARG:HD2	2.21	0.41
1:C:399:LEU:O	1:C:452:LYS:NZ	2.52	0.41
1:A:549:GLU:OE2	4:A:942:HOH:O	2.22	0.41
1:A:502:GLN:H	2:B:0:ALA:C	2.25	0.41
1:C:444:ASN:OD1	1:C:477:LEU:HD21	2.21	0.40
1:A:483:THR:HG21	1:A:485:ASN:OD1	2.17	0.40
1:A:462:ASP:HA	1:A:491:ALA:HB2	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 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	377/388 (97%)	366 (97%)	10 (3%)	1 (0%)	46	63
1	C	378/388 (97%)	366 (97%)	12 (3%)	0	100	100
2	B	2/5 (40%)	2 (100%)	0	0	100	100
2	D	2/5 (40%)	2 (100%)	0	0	100	100
All	All	759/786 (97%)	736 (97%)	22 (3%)	1 (0%)	56	74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	710	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	336/345 (97%)	316 (94%)	20 (6%)	24	37
1	C	337/345 (98%)	316 (94%)	21 (6%)	23	35
2	B	3/3 (100%)	3 (100%)	0	100	100
2	D	3/3 (100%)	3 (100%)	0	100	100
All	All	679/696 (98%)	638 (94%)	41 (6%)	24	37

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	338	MET
1	A	341	LYS
1	A	351	MET
1	A	363	LEU
1	A	403	ASP
1	A	472	ASP
1	A	483	THR
1	A	485	ASN
1	A	541	LEU
1	A	553	SER
1	A	564	ILE
1	A	621	LYS
1	A	638	ASP
1	A	658	LEU
1	A	660	SER
1	A	680	GLU
1	A	706	ASN
1	A	709	LYS
1	A	711	LEU
1	A	716	ASP
1	C	351	MET
1	C	353	TYR
1	C	401	LEU
1	C	447	LYS
1	C	467	ARG
1	C	473	THR
1	C	483	THR
1	C	485	ASN
1	C	486	ILE
1	C	506	LEU
1	C	565	GLN
1	C	573	SER
1	C	578	LEU
1	C	585	GLU
1	C	662	ASP
1	C	665	LYS
1	C	668	LEU
1	C	679	LYS
1	C	696	GLU
1	C	709	LYS
1	C	712	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	450	GLN
1	A	485	ASN
1	A	565	GLN
1	A	599	GLN
1	A	681	HIS
1	C	485	ASN
1	C	577	ASN
1	C	584	HIS
1	C	601	GLN
1	C	681	HIS
1	C	691	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	4AR	B	4	1,2	7,11,23	1.89	1 (14%)	6,13,30	7.34	2 (33%)
2	4AR	D	4	1,2	7,11,23	5.64	1 (14%)	6,13,30	7.41	2 (33%)

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	4AR	B	4	1,2	-	0/8/11/20	0/0/0/1
2	4AR	D	4	1,2	-	0/8/11/20	0/0/0/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	4	4AR	C7-C2	4.90	1.63	1.49
2	D	4	4AR	C7-C2	14.89	1.92	1.49

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	4	4AR	O1-C2-C7	-16.92	87.94	121.01
2	D	4	4AR	O1-C2-C7	-13.78	94.07	121.01
2	D	4	4AR	C7-C2-C1	-11.77	85.96	116.70
2	B	4	4AR	C7-C2-C1	-6.08	100.80	116.70

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	4	4AR	3	0
2	D	4	4AR	7	0

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

10 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	$\# Z > 2$
3	ACT	A	801	-	1,3,3	0.52	0	0,3,3	0.00	-
3	ACT	A	802	-	1,3,3	0.75	0	0,3,3	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ACT	A	803	-	1,3,3	0.61	0	0,3,3	0.00	-
3	ACT	A	804	-	1,3,3	0.39	0	0,3,3	0.00	-
3	ACT	A	805	-	1,3,3	0.76	0	0,3,3	0.00	-
3	ACT	C	801	-	1,3,3	1.28	0	0,3,3	0.00	-
3	ACT	C	802	-	1,3,3	0.04	0	0,3,3	0.00	-
3	ACT	C	803	-	1,3,3	1.19	0	0,3,3	0.00	-
3	ACT	C	804	-	1,3,3	0.40	0	0,3,3	0.00	-
3	ACT	C	805	-	1,3,3	1.14	0	0,3,3	0.00	-

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
3	ACT	A	801	-	-	0/0/0/0	0/0/0/0
3	ACT	A	802	-	-	0/0/0/0	0/0/0/0
3	ACT	A	803	-	-	0/0/0/0	0/0/0/0
3	ACT	A	804	-	-	0/0/0/0	0/0/0/0
3	ACT	A	805	-	-	0/0/0/0	0/0/0/0
3	ACT	C	801	-	-	0/0/0/0	0/0/0/0
3	ACT	C	802	-	-	0/0/0/0	0/0/0/0
3	ACT	C	803	-	-	0/0/0/0	0/0/0/0
3	ACT	C	804	-	-	0/0/0/0	0/0/0/0
3	ACT	C	805	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	801	ACT	8	0
3	A	803	ACT	2	0
3	A	804	ACT	2	0
3	C	803	ACT	1	0
3	C	805	ACT	1	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.

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	379/388 (97%)	0.19	19 (5%) 32 33	42, 66, 124, 155	29 (7%)
1	C	379/388 (97%)	0.22	21 (5%) 29 29	42, 66, 122, 151	27 (7%)
2	B	4/5 (80%)	1.13	0 100 100	61, 69, 75, 84	0
2	D	4/5 (80%)	0.80	0 100 100	64, 78, 84, 87	0
All	All	766/786 (97%)	0.21	40 (5%) 31 31	42, 66, 122, 155	56 (7%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	504	SER	6.4
1	A	697	ASP	4.5
1	C	580	TRP	4.5
1	C	663	LEU	4.1
1	A	505	GLY	4.0
1	C	583	ALA	3.9
1	A	710	PRO	3.9
1	A	695	LEU	3.6
1	A	586	LEU	3.5
1	C	573	SER	3.5
1	A	620	TYR	3.1
1	C	659	VAL	3.0
1	C	584	HIS	3.0
1	C	664	PRO	3.0
1	A	708	GLY	2.9
1	A	578	LEU	2.9
1	A	600	ILE	2.8
1	C	617	SER	2.8
1	C	699	VAL	2.6
1	C	567	THR	2.6
1	A	580	TRP	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	619	VAL	2.6
1	C	576	ARG	2.6
1	C	575	VAL	2.5
1	A	591	CYS	2.5
1	C	665	LYS	2.5
1	A	506	LEU	2.5
1	A	643	PRO	2.4
1	A	658	LEU	2.4
1	C	507	ALA	2.3
1	C	712	ILE	2.3
1	A	657	TYR	2.3
1	A	667	CYS	2.2
1	C	717	MET	2.2
1	C	570	SER	2.1
1	A	401	LEU	2.1
1	C	658	LEU	2.1
1	C	506	LEU	2.1
1	C	569	TYR	2.1
1	A	583	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	4AR	B	4	12/23	0.93	0.17	-	48,51,59,124	0
2	4AR	D	4	12/23	0.91	0.16	-	49,53,56,98	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	ACT	C	801	4/4	0.74	0.34	25.16	88,89,90,92	0
3	ACT	A	803	4/4	0.77	0.38	18.25	82,83,84,86	0
3	ACT	A	805	4/4	0.63	0.40	10.34	109,110,111,112	0
3	ACT	A	804	4/4	0.90	0.33	7.39	81,82,83,86	0
3	ACT	A	801	4/4	0.87	0.29	5.38	77,80,80,82	0
3	ACT	C	805	4/4	0.15	0.37	3.56	122,125,126,126	0
3	ACT	C	803	4/4	0.78	0.33	2.33	98,100,100,100	0
3	ACT	C	804	4/4	0.91	0.19	-	97,97,98,98	0
3	ACT	C	802	4/4	0.90	0.28	-	94,95,95,96	0
3	ACT	A	802	4/4	0.87	0.30	-	106,107,108,109	0

6.5 Other polymers [i](#)

There are no such residues in this entry.