



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:56 PM GMT

PDB ID : 5BV1  
Title : Crystal Structure of a Vps33-Vps16 Complex from Chaetomium thermophilum  
Authors : Baker, R.W.; Jeffrey, P.D.; Hughson, F.M.  
Deposited on : 2015-06-04  
Resolution : 2.90 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
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) : rb-20026982

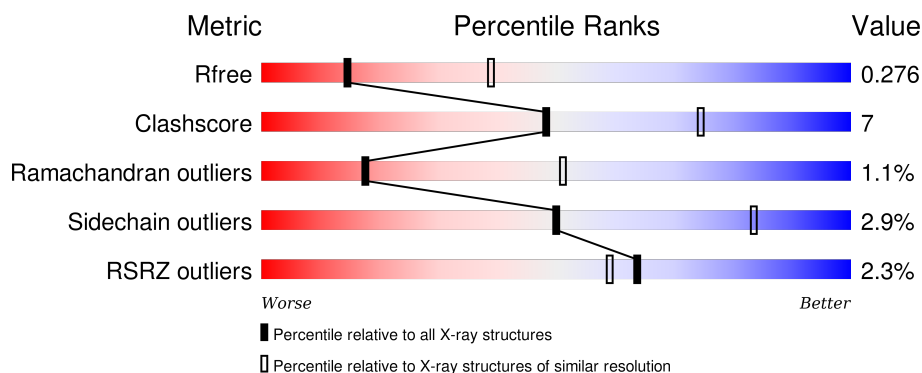
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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  $\geq 3$ , 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	669	<div> <div>2%</div> <div>71% 18% • 10%</div> </div>
1	C	669	<div> <div>%</div> <div>72% 16% • 10%</div> </div>
2	B	333	<div> <div>2%</div> <div>67% 21% • 11%</div> </div>
2	D	333	<div> <div>5%</div> <div>73% 15% • 11%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14259 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 VPS33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	601	Total	C	N	O	S	0	0	0
			4752	3007	836	898	11			
1	C	602	Total	C	N	O	S	0	0	0
			4759	3012	837	899	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP G0SCM5
C	-1	GLY	-	expression tag	UNP G0SCM5

- Molecule 2 is a protein called Putative vacuolar protein sorting-associated protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	297	Total	C	N	O	S	0	0	0
			2365	1487	423	446	9			
2	D	297	Total	C	N	O	S	0	0	0
			2365	1487	423	446	9			

There are 44 discrepancies between the modelled and reference sequences:

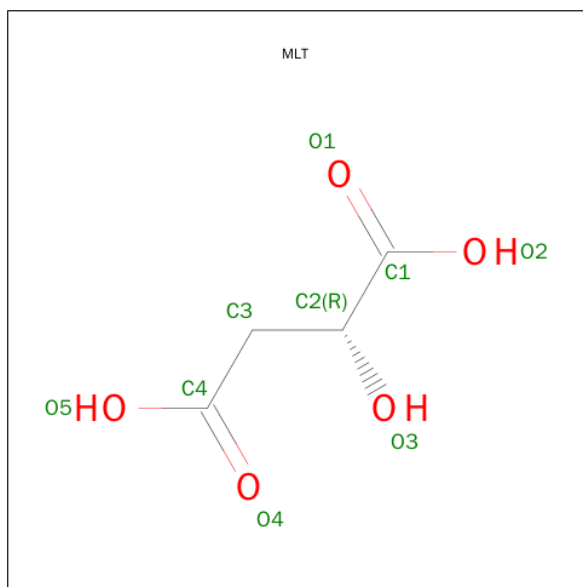
Chain	Residue	Modelled	Actual	Comment	Reference
B	502	MET	-	initiating methionine	UNP G0S6M7
B	503	GLY	-	expression tag	UNP G0S6M7
B	504	SER	-	expression tag	UNP G0S6M7
B	672	ARG	-	insertion	UNP G0S6M7
B	673	MET	-	insertion	UNP G0S6M7
B	674	GLN	-	insertion	UNP G0S6M7
B	675	GLU	-	insertion	UNP G0S6M7
B	676	THR	-	insertion	UNP G0S6M7
B	677	PHE	-	insertion	UNP G0S6M7
B	678	GLU	-	insertion	UNP G0S6M7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	679	ARG	-	insertion	UNP G0S6M7
B	680	ASP	-	insertion	UNP G0S6M7
B	681	LEU	-	insertion	UNP G0S6M7
B	682	THR	-	insertion	UNP G0S6M7
B	683	ASP	-	insertion	UNP G0S6M7
B	684	SER	-	insertion	UNP G0S6M7
B	685	PHE	-	insertion	UNP G0S6M7
B	686	VAL	-	insertion	UNP G0S6M7
B	687	GLY	-	insertion	UNP G0S6M7
B	688	LEU	-	insertion	UNP G0S6M7
B	689	SER	-	insertion	UNP G0S6M7
B	690	VAL	LEU	engineered mutation	UNP G0S6M7
D	502	MET	-	initiating methionine	UNP G0S6M7
D	503	GLY	-	expression tag	UNP G0S6M7
D	504	SER	-	expression tag	UNP G0S6M7
D	672	ARG	-	insertion	UNP G0S6M7
D	673	MET	-	insertion	UNP G0S6M7
D	674	GLN	-	insertion	UNP G0S6M7
D	675	GLU	-	insertion	UNP G0S6M7
D	676	THR	-	insertion	UNP G0S6M7
D	677	PHE	-	insertion	UNP G0S6M7
D	678	GLU	-	insertion	UNP G0S6M7
D	679	ARG	-	insertion	UNP G0S6M7
D	680	ASP	-	insertion	UNP G0S6M7
D	681	LEU	-	insertion	UNP G0S6M7
D	682	THR	-	insertion	UNP G0S6M7
D	683	ASP	-	insertion	UNP G0S6M7
D	684	SER	-	insertion	UNP G0S6M7
D	685	PHE	-	insertion	UNP G0S6M7
D	686	VAL	-	insertion	UNP G0S6M7
D	687	GLY	-	insertion	UNP G0S6M7
D	688	LEU	-	insertion	UNP G0S6M7
D	689	SER	-	insertion	UNP G0S6M7
D	690	VAL	LEU	engineered mutation	UNP G0S6M7

- Molecule 3 is D-MALATE (three-letter code: MLT) (formula: C<sub>4</sub>H<sub>6</sub>O<sub>5</sub>).

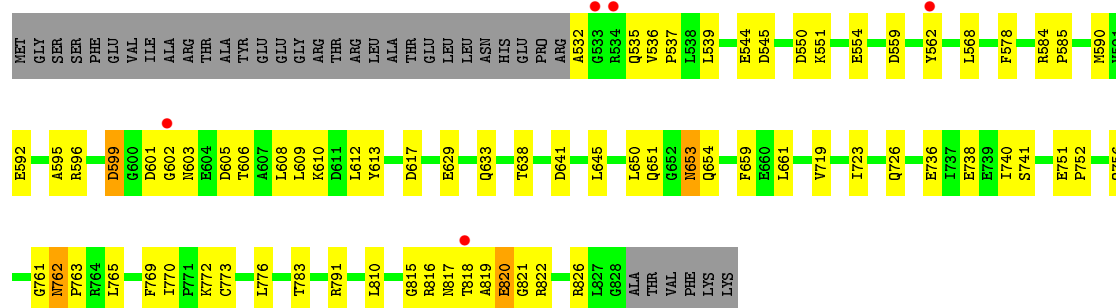


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			9	4	5		
3	D	1	Total	C	O	0	0
			9	4	5		

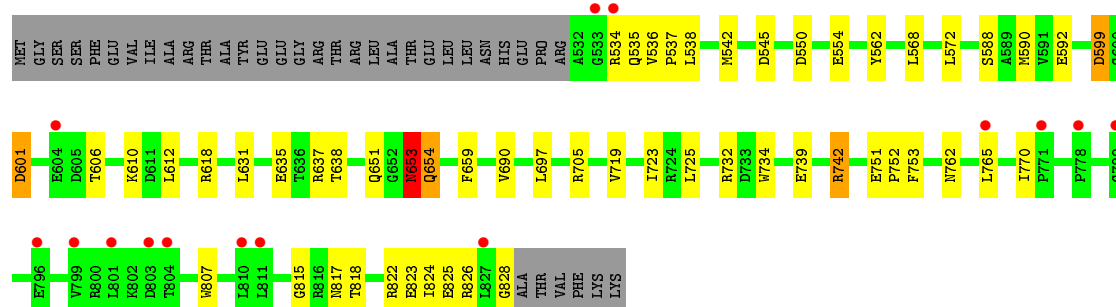
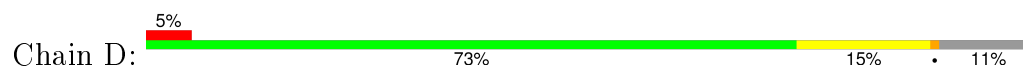




- Molecule 2: Putative vacuolar protein sorting-associated protein



- Molecule 2: Putative vacuolar protein sorting-associated protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.28Å 159.90Å 97.39Å 90.00° 90.21° 90.00°	Depositor
Resolution (Å)	61.59 – 2.90 83.18 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.9 (61.59-2.90) 98.9 (83.18-2.90)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.15 (at 2.91Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.211 , 0.276 0.213 , 0.276	Depositor DCC
$R_{free}$ test set	2699 reflections (5.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	65.3	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 32.8	EDS
Estimated twinning fraction	0.155 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 53053 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	14259	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MLT

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	0.34	0/4827	0.54	0/6512
1	C	0.33	0/4834	0.54	0/6522
2	B	0.32	0/2404	0.50	0/3242
2	D	0.32	0/2404	0.54	0/3242
All	All	0.33	0/14469	0.54	0/19518

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	C	0	2
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	340	ASN	Peptide
1	A	341	THR	Peptide
1	C	340	ASN	Peptide
1	C	341	THR	Peptide

## 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	4752	0	4794	76	0
1	C	4759	0	4803	68	1
2	B	2365	0	2368	44	0
2	D	2365	0	2368	33	1
3	B	9	0	4	1	0
3	D	9	0	4	1	0
All	All	14259	0	14341	210	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 7.

All (210) 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:A:102:GLN:HE21	1:A:103:THR:H	1.20	0.89
1:C:153:LEU:HB2	1:C:156:SER:HB3	1.65	0.77
1:A:31:LYS:NZ	1:A:55:TYR:O	2.18	0.75
1:A:338:ARG:O	1:A:341:THR:OG1	2.02	0.75
1:A:231:GLU:OE2	1:A:560:GLN:NE2	2.22	0.73
1:C:181:LYS:HE2	1:C:210:LEU:HD11	1.70	0.72
2:D:545:ASP:HB3	2:D:568:LEU:HD21	1.71	0.71
1:A:95:LYS:HA	1:A:98:GLN:HB2	1.72	0.71
1:A:70:THR:OG1	1:A:102:GLN:OE1	2.08	0.71
1:A:382:ILE:HD13	1:A:414:THR:HG22	1.75	0.69
2:D:653:ASN:ND2	3:D:901:MLT:O4	2.25	0.69
1:C:505:GLU:OE1	1:C:505:GLU:N	2.27	0.67
1:C:500:ARG:NH2	1:C:505:GLU:OE2	2.27	0.67
2:B:606:THR:HG22	2:B:610:LYS:HE3	1.77	0.66
2:B:820:GLU:C	2:B:822:ARG:H	2.00	0.66
1:A:153:LEU:HB2	1:A:156:SER:HB3	1.80	0.64
1:C:68:THR:OG1	1:C:93:GLN:NE2	2.32	0.63
1:C:499:LEU:O	1:C:527:ARG:HD2	1.99	0.62
1:A:505:GLU:N	1:A:505:GLU:OE1	2.32	0.62
1:C:103:THR:OG1	1:C:104:SER:N	2.33	0.62
1:A:181:LYS:HE2	1:A:210:LEU:HD11	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:822:ARG:O	2:B:826:ARG:N	2.30	0.61
1:C:160:LEU:HD11	1:C:169:VAL:HG23	1.81	0.61
1:A:479:MET:SD	1:A:570:LYS:HD3	2.42	0.60
1:A:607:VAL:HG22	1:A:636:CYS:HB2	1.84	0.59
1:A:137:LEU:HD22	1:A:139:LEU:HG	1.85	0.59
1:C:160:LEU:HD13	1:C:168:PRO:HB2	1.84	0.59
2:D:822:ARG:O	2:D:826:ARG:N	2.30	0.59
1:A:438:LEU:HD22	1:A:468:SER:HB2	1.84	0.58
2:D:823:GLU:HA	2:D:826:ARG:HB2	1.86	0.57
1:A:432:GLY:O	1:A:469:LYS:NZ	2.35	0.56
1:C:507:ASN:OD1	1:C:510:ASP:N	2.38	0.56
1:A:499:LEU:O	1:A:527:ARG:HD2	2.06	0.56
1:A:500:ARG:O	1:A:527:ARG:NH1	2.39	0.56
1:A:53:ARG:HG2	1:A:57:VAL:O	2.06	0.56
1:A:85:ARG:NH1	2:B:601:ASP:OD2	2.40	0.55
1:A:73:ARG:NH1	1:A:103:THR:OG1	2.40	0.54
1:C:65:ASN:HB3	1:C:67:ASN:OD1	2.07	0.54
1:C:338:ARG:O	1:C:341:THR:OG1	2.26	0.54
1:A:160:LEU:HD11	1:A:169:VAL:HG23	1.90	0.53
1:A:330:ARG:NH1	1:A:360:GLU:OE2	2.42	0.52
2:D:550:ASP:O	2:D:554:GLU:HG3	2.10	0.52
1:A:507:ASN:OD1	1:A:510:ASP:N	2.30	0.52
2:B:820:GLU:OE1	2:B:826:ARG:NH2	2.42	0.52
1:A:85:ARG:HD2	2:B:601:ASP:CG	2.29	0.52
2:D:606:THR:HG22	2:D:610:LYS:HE3	1.92	0.52
1:A:456:THR:HG23	1:A:651:ILE:HA	1.92	0.52
2:D:534:ARG:HA	2:D:537:PRO:HD2	1.92	0.52
1:C:456:THR:HG23	1:C:651:ILE:HA	1.92	0.51
1:A:301:GLN:NE2	1:A:303:ASP:OD2	2.37	0.51
1:C:95:LYS:HA	1:C:98:GLN:HB2	1.93	0.51
1:C:494:TYR:CZ	1:C:498:GLN:HG3	2.46	0.51
2:D:568:LEU:HD22	2:D:572:LEU:HD11	1.93	0.51
1:A:102:GLN:NE2	1:A:103:THR:H	2.00	0.51
1:A:526:ILE:HD12	1:A:620:LEU:HD13	1.92	0.51
2:B:545:ASP:HB3	2:B:568:LEU:HD21	1.92	0.51
2:B:641:ASP:O	2:B:645:LEU:HG	2.11	0.51
2:B:592:GLU:OE2	2:B:613:TYR:OH	2.25	0.50
1:C:607:VAL:HG22	1:C:636:CYS:HB2	1.92	0.50
2:B:596:ARG:HG2	2:B:609:LEU:HD11	1.94	0.50
1:C:88:HIS:HB3	2:D:599:ASP:O	2.11	0.50
1:A:91:ALA:HB2	1:A:123:LEU:HD23	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:ARG:NH1	1:A:58:ASP:O	2.44	0.50
1:C:161:TYR:CE1	1:C:239:GLU:HB3	2.47	0.49
1:C:33:LEU:HD12	1:C:76:VAL:HB	1.94	0.49
2:B:726:GLN:HG2	2:B:756:GLN:NE2	2.27	0.49
1:C:121:LYS:HA	1:C:124:GLU:HG2	1.95	0.49
1:C:191:GLY:HA3	1:C:577:PHE:CZ	2.48	0.49
2:B:815:GLY:O	2:B:817:ASN:N	2.44	0.49
1:C:268:ASP:HB3	1:C:297:LYS:HE2	1.94	0.49
1:C:84:VAL:HB	1:C:118:PHE:CZ	2.47	0.49
2:B:820:GLU:C	2:B:822:ARG:N	2.66	0.48
1:A:517:VAL:HG23	1:A:518:TYR:CD2	2.48	0.48
1:A:84:VAL:HB	1:A:118:PHE:CZ	2.47	0.48
2:D:807:TRP:CH2	2:D:828:GLY:HA3	2.48	0.48
1:A:494:TYR:CZ	1:A:498:GLN:HG3	2.48	0.48
1:C:386:LEU:HD23	1:C:422:LEU:HD11	1.94	0.48
1:C:36:GLU:OE2	1:C:80:ARG:HG3	2.12	0.48
1:A:191:GLY:HA3	1:A:577:PHE:CZ	2.49	0.48
1:C:338:ARG:HB3	1:C:350:PHE:CD1	2.48	0.48
1:A:30:LYS:HB3	1:A:30:LYS:HE2	1.54	0.48
2:B:536:VAL:HB	2:B:537:PRO:HD3	1.96	0.48
1:A:174:ARG:HA	1:A:206:MET:HE1	1.96	0.47
1:A:160:LEU:HD23	1:A:161:TYR:CZ	2.49	0.47
2:D:807:TRP:HZ2	2:D:825:GLU:HA	1.80	0.47
2:D:762:ASN:OD1	2:D:765:LEU:N	2.36	0.47
1:C:161:TYR:OH	1:C:237:ASP:OD1	2.30	0.47
1:C:269:VAL:O	1:C:298:ARG:N	2.33	0.47
1:C:182:LYS:HG2	1:C:183:HIS:CE1	2.50	0.47
1:A:408:GLU:O	1:A:412:ARG:HG3	2.14	0.47
1:A:97:ILE:O	1:A:101:SER:HB3	2.15	0.47
2:B:773:CYS:HB3	2:B:776:LEU:HD13	1.96	0.47
1:A:354:LEU:N	1:A:355:PRO:HD2	2.29	0.47
2:D:536:VAL:HB	2:D:537:PRO:HD3	1.96	0.47
2:D:592:GLU:OE2	2:D:618:ARG:NH2	2.48	0.47
1:A:103:THR:OG1	1:A:104:SER:N	2.46	0.47
2:D:653:ASN:HB3	2:D:654:GLN:H	1.50	0.47
1:C:500:ARG:O	1:C:527:ARG:NH1	2.48	0.47
1:C:654:ALA:HB3	2:D:631:LEU:HD13	1.97	0.46
2:D:742:ARG:HE	2:D:742:ARG:HB2	1.56	0.46
1:C:186:PHE:O	1:C:581:GLN:NE2	2.37	0.46
2:D:732:ARG:HD2	2:D:734:TRP:CZ2	2.50	0.46
1:C:252:TYR:O	1:C:256:ILE:HG13	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:539:LEU:O	2:B:544:GLU:HB2	2.15	0.46
1:C:338:ARG:HB3	1:C:350:PHE:CE1	2.50	0.46
1:C:174:ARG:HA	1:C:206:MET:HE1	1.97	0.46
1:A:192:LYS:HE2	1:A:572:ALA:HB3	1.97	0.46
2:B:819:ALA:O	2:B:821:GLY:N	2.49	0.46
1:A:354:LEU:H	1:A:355:PRO:HD2	1.81	0.46
1:A:116:THR:HB	2:B:559:ASP:OD1	2.15	0.46
1:A:386:LEU:HD23	1:A:422:LEU:HD11	1.97	0.46
1:C:655:THR:O	1:C:656:PHE:HB2	2.15	0.46
1:A:174:ARG:HG2	1:A:206:MET:HE3	1.98	0.46
1:C:7:PHE:CZ	1:C:149:LEU:HD11	2.51	0.46
1:C:188:ARG:NH1	1:C:578:ASP:OD2	2.48	0.46
2:B:606:THR:CG2	2:B:610:LYS:HE3	2.46	0.45
1:C:85:ARG:HG3	2:D:601:ASP:N	2.31	0.45
1:A:655:THR:O	1:A:656:PHE:HB2	2.15	0.45
1:C:90:ILE:HD12	1:C:111:TRP:CH2	2.52	0.45
1:A:114:ARG:HH22	2:B:559:ASP:HB2	1.81	0.45
1:A:114:ARG:NH2	2:B:559:ASP:HB2	2.31	0.45
2:B:751:GLU:HB2	2:B:752:PRO:HD3	1.98	0.45
1:A:500:ARG:NH2	1:A:505:GLU:OE2	2.50	0.45
2:D:807:TRP:NE1	2:D:824:ILE:O	2.43	0.44
1:A:63:LEU:HD12	1:A:93:GLN:HG3	1.98	0.44
1:A:33:LEU:HD22	1:A:57:VAL:HG11	1.99	0.44
1:A:335:TYR:HE1	1:A:354:LEU:HD11	1.83	0.44
1:A:145:GLU:OE2	1:A:304:GLY:N	2.43	0.44
1:C:386:LEU:O	1:C:390:GLN:HG3	2.17	0.44
2:B:762:ASN:HB2	2:B:765:LEU:H	1.82	0.44
2:B:791:ARG:HG3	2:B:810:LEU:HD22	2.00	0.44
2:B:617:ASP:O	3:B:901:MLT:H31	2.17	0.44
2:B:532:ALA:HA	2:B:535:GLN:HB3	1.98	0.44
1:C:249:GLN:HB3	1:C:615:THR:HG23	2.00	0.44
2:B:815:GLY:C	2:B:817:ASN:H	2.20	0.44
2:D:751:GLU:HB2	2:D:752:PRO:HD3	1.98	0.44
1:C:442:ARG:HB3	2:D:659:PHE:CZ	2.52	0.44
2:D:534:ARG:CA	2:D:537:PRO:HD2	2.48	0.44
1:C:438:LEU:HG	1:C:442:ARG:HD2	1.99	0.44
2:B:650:LEU:HB2	2:B:661:LEU:HD21	2.00	0.44
2:B:595:ALA:O	2:B:605:ASP:N	2.41	0.44
1:A:102:GLN:NE2	1:A:103:THR:HG22	2.32	0.44
1:C:364:LEU:O	1:C:368:SER:OG	2.23	0.44
1:C:84:VAL:HB	1:C:118:PHE:CE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:ASP:HB3	1:A:297:LYS:HE2	1.99	0.43
1:C:266:GLN:OE1	1:C:299:LYS:HE2	2.18	0.43
1:A:337:SER:O	1:A:339:HIS:N	2.51	0.43
2:D:697:LEU:HD13	2:D:705:ARG:HB3	2.01	0.43
1:C:526:ILE:HD12	1:C:620:LEU:HG	1.99	0.43
1:A:108:HIS:ND1	1:A:133:ASN:HB2	2.34	0.43
1:A:112:VAL:HA	1:A:113:PRO:HA	1.75	0.43
1:A:331:LEU:HA	1:A:331:LEU:HD23	1.88	0.43
1:C:31:LYS:NZ	1:C:55:TYR:O	2.52	0.43
1:A:495:LEU:HD22	1:A:499:LEU:HD12	2.01	0.43
1:A:526:ILE:O	1:A:530:GLN:HG3	2.19	0.43
2:D:588:SER:O	2:D:592:GLU:HB2	2.19	0.43
1:A:27:VAL:HB	1:A:31:LYS:HE2	2.01	0.42
1:C:494:TYR:CE2	1:C:498:GLN:HG3	2.54	0.42
1:C:538:LEU:HD12	1:C:538:LEU:HA	1.90	0.42
1:A:102:GLN:HE21	1:A:103:THR:HG22	1.83	0.42
1:A:303:ASP:OD1	1:A:305:SER:OG	2.19	0.42
2:D:562:TYR:HB2	2:D:590:MET:CE	2.50	0.42
1:A:641:ILE:HG13	1:A:642:ASN:N	2.34	0.42
1:C:7:PHE:HA	1:C:183:HIS:NE2	2.35	0.42
1:C:151:LEU:HD11	1:C:172:LEU:HA	2.01	0.42
2:B:738:GLU:O	2:B:741:SER:OG	2.27	0.42
2:B:629:GLU:O	2:B:633:GLN:HG2	2.18	0.42
2:D:725:LEU:HD13	2:D:753:PHE:CD2	2.54	0.42
2:D:719:VAL:O	2:D:723:ILE:HG13	2.19	0.42
1:A:319:ILE:HD13	1:A:509:GLN:HA	2.00	0.42
2:B:736:GLU:O	2:B:740:ILE:HG13	2.20	0.42
1:C:53:ARG:HG2	1:C:57:VAL:O	2.20	0.42
1:C:186:PHE:O	1:C:581:GLN:HB2	2.20	0.42
1:A:188:ARG:NH1	1:A:578:ASP:OD2	2.53	0.42
2:B:578:PHE:HZ	2:B:608:LEU:HD12	1.84	0.42
1:C:495:LEU:HD22	1:C:499:LEU:CD1	2.50	0.41
1:A:442:ARG:HB3	2:B:659:PHE:CZ	2.55	0.41
1:C:517:VAL:HG23	1:C:518:TYR:CD2	2.55	0.41
1:C:426:TYR:HD1	1:C:426:TYR:C	2.23	0.41
1:C:499:LEU:HD22	1:C:527:ARG:HB3	2.03	0.41
1:C:174:ARG:HG2	1:C:206:MET:HE3	2.02	0.41
2:B:751:GLU:CD	2:B:751:GLU:H	2.23	0.41
2:D:635:GLU:HG3	2:D:637:ARG:H	1.85	0.41
1:C:354:LEU:N	1:C:355:PRO:HD2	2.35	0.41
2:B:578:PHE:CD1	2:B:612:LEU:HD13	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:GLN:HE22	1:C:183:HIS:CE1	2.38	0.41
1:C:655:THR:HG23	1:C:657:GLU:H	1.86	0.41
2:B:584:ARG:HA	2:B:585:PRO:HD2	1.97	0.41
2:B:550:ASP:O	2:B:554:GLU:HG3	2.21	0.41
2:B:769:PHE:HA	2:B:772:LYS:HD3	2.02	0.41
2:D:538:LEU:O	2:D:542:MET:HG3	2.21	0.41
1:A:424:CYS:SG	1:A:460:LEU:HD13	2.61	0.41
1:A:374:ILE:O	1:A:378:THR:HG23	2.21	0.41
2:B:551:LYS:HA	2:B:551:LYS:HD3	1.88	0.41
2:B:599:ASP:OD2	2:B:603:ASN:HB2	2.20	0.41
1:A:160:LEU:HD13	1:A:168:PRO:HB2	2.02	0.41
1:C:447:GLN:HG2	2:D:690:VAL:HG21	2.02	0.41
1:A:24:LEU:HD23	1:A:24:LEU:HA	1.93	0.41
1:C:201:ASP:O	1:C:205:ARG:HG3	2.21	0.41
1:A:601:LYS:HB2	1:A:630:ARG:C	2.42	0.41
2:D:815:GLY:O	2:D:817:ASN:N	2.53	0.41
1:C:426:TYR:C	1:C:426:TYR:CD1	2.94	0.40
1:A:6:GLY:HA2	1:A:229:THR:HG21	2.03	0.40
2:D:807:TRP:CZ2	2:D:825:GLU:HA	2.57	0.40
1:C:356:GLY:O	1:C:360:GLU:HB2	2.21	0.40
1:C:145:GLU:OE2	1:C:304:GLY:N	2.49	0.40
2:B:761:GLY:O	2:B:763:PRO:HD3	2.21	0.40
1:A:537:TYR:O	1:A:540:SER:HB3	2.22	0.40
2:B:562:TYR:HB2	2:B:590:MET:CE	2.51	0.40
2:B:719:VAL:O	2:B:723:ILE:HG13	2.22	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:C:342:LYS:NZ	2:D:535:GLN:OE1[1_455]	2.19	0.01

## 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	593/669 (89%)	552 (93%)	35 (6%)	6 (1%)	19	54
1	C	594/669 (89%)	548 (92%)	40 (7%)	6 (1%)	19	54
2	B	295/333 (89%)	267 (90%)	23 (8%)	5 (2%)	11	38
2	D	295/333 (89%)	266 (90%)	26 (9%)	3 (1%)	19	54
All	All	1777/2004 (89%)	1633 (92%)	124 (7%)	20 (1%)	17	51

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	338	ARG
1	A	482	SER
2	B	653	ASN
2	B	654	GLN
1	C	338	ARG
2	D	601	ASP
2	D	653	ASN
2	D	654	GLN
1	A	101	SER
2	B	602	GLY
1	C	271	ALA
1	C	337	SER
1	C	342	LYS
1	A	102	GLN
1	A	219	SER
2	B	820	GLU
1	C	219	SER
2	B	816	ARG
1	C	102	GLN
1	A	486	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	515/564 (91%)	504 (98%)	11 (2%)	61	88
1	C	516/564 (92%)	500 (97%)	16 (3%)	47	82
2	B	248/280 (89%)	240 (97%)	8 (3%)	46	81
2	D	248/280 (89%)	239 (96%)	9 (4%)	42	78
All	All	1527/1688 (90%)	1483 (97%)	44 (3%)	50	83

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	GLN
1	A	131	ASP
1	A	135	SER
1	A	137	LEU
1	A	207	ARG
1	A	324	LEU
1	A	330	ARG
1	A	426	TYR
1	A	542	THR
1	A	603	THR
1	A	620	LEU
2	B	599	ASP
2	B	638	THR
2	B	651	GLN
2	B	653	ASN
2	B	762	ASN
2	B	770	ILE
2	B	783	THR
2	B	818	THR
1	C	11	GLN
1	C	102	GLN
1	C	114	ARG
1	C	131	ASP
1	C	135	SER
1	C	207	ARG
1	C	225	SER
1	C	272	VAL
1	C	299	LYS
1	C	330	ARG
1	C	338	ARG
1	C	426	TYR
1	C	454	LEU
1	C	499	LEU

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Mol	Chain	Res	Type
1	C	533	LEU
1	C	603	THR
2	D	599	ASP
2	D	612	LEU
2	D	638	THR
2	D	651	GLN
2	D	653	ASN
2	D	739	GLU
2	D	742	ARG
2	D	770	ILE
2	D	818	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	183	HIS

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	MLT	B	901	-	1,8,8	0.53	0	2,10,10	4.72	1 (50%)
3	MLT	D	901	-	1,8,8	0.69	0	2,10,10	2.41	1 (50%)

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	MLT	B	901	-	-	0/2/8/8	0/0/0/0
3	MLT	D	901	-	-	0/2/8/8	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	B	901	MLT	C4-C3-C2	-6.68	107.24	114.33
3	D	901	MLT	C4-C3-C2	-3.09	111.06	114.33

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	901	MLT	1	0
3	D	901	MLT	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	601/669 (89%)	0.16	14 (2%) 64 59	30, 54, 120, 170	0
1	C	602/669 (89%)	0.15	8 (1%) 79 78	30, 55, 113, 152	0
2	B	297/333 (89%)	0.12	5 (1%) 73 70	48, 65, 106, 134	0
2	D	297/333 (89%)	0.44	15 (5%) 32 25	42, 72, 141, 165	0
All	All	1797/2004 (89%)	0.20	42 (2%) 64 59	30, 60, 123, 170	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	75	ILE	6.0
2	D	533	GLY	5.1
1	C	221	ARG	4.4
1	C	483	SER	4.1
1	A	219	SER	4.1
1	A	482	SER	4.1
2	B	533	GLY	4.1
1	A	221	ARG	4.0
1	C	342	LYS	3.9
1	A	342	LYS	3.7
2	D	799	VAL	3.6
2	D	803	ASP	3.5
2	D	827	LEU	3.4
2	B	534	ARG	3.2
2	D	810	LEU	3.0
2	D	604	GLU	3.0
1	A	343	THR	2.9
1	C	484	GLY	2.9
2	D	804	THR	2.8
2	B	562	TYR	2.8
2	D	811	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	212	ALA	2.6
2	D	789	GLY	2.6
2	D	771	PRO	2.6
2	D	534	ARG	2.5
1	A	94	ILE	2.5
1	C	340	ASN	2.4
1	A	99	ARG	2.4
1	C	127	GLY	2.4
2	B	602	GLY	2.4
1	A	128	VAL	2.4
2	B	818	THR	2.3
1	A	344	THR	2.3
1	A	77	PHE	2.3
2	D	765	LEU	2.2
1	A	483	SER	2.2
2	D	796	GLU	2.1
2	D	801	LEU	2.1
1	A	122	VAL	2.1
1	C	118	PHE	2.1
1	A	220	ASP	2.1
2	D	778	PRO	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	MLT	B	901	9/9	0.88	0.25	0.57	58,70,89,99	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	MLT	D	901	9/9	0.88	0.17	-0.85	72,75,89,95	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.