



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:55 AM GMT

PDB ID : 3K8P
Title : Structural basis for vesicle tethering by the Dsl1 complex
Authors : Ren, Y.; Jeffrey, P.D.; Hughson, F.M.
Deposited on : 2009-10-14
Resolution : 2.60 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

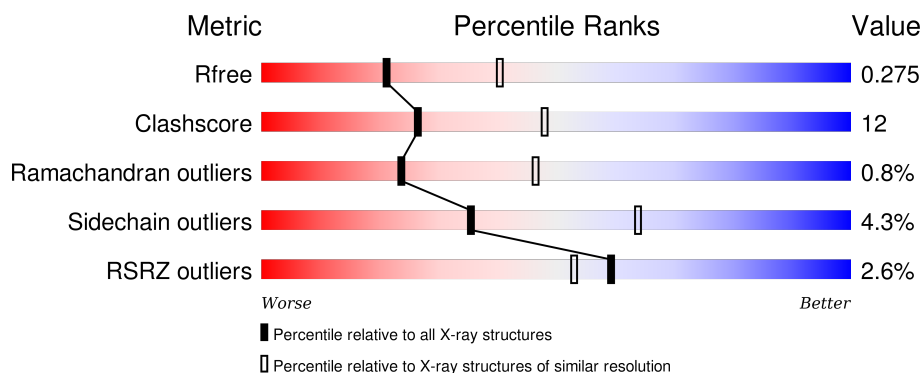
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	C	357	<div> <div>2%</div> <div>62%</div> <div>19%</div> <div>•</div> <div>17%</div> </div>
2	D	709	<div> <div>2%</div> <div>59%</div> <div>22%</div> <div>•</div> <div>16%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7376 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 Dsl1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	C	295	Total	C	N	O	S	Se	0	0	0
			2464	1594	404	458	2	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	330	GLY	-	EXPRESSION TAG	UNP Q6CUS2
C	331	SER	-	EXPRESSION TAG	UNP Q6CUS2

- Molecule 2 is a protein called Protein transport protein SEC39.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	D	594	Total	C	N	O	S	Se	0	0	0
			4909	3182	775	931	7	14			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	3	Total	O	0	0
			3	3		

Y631	D632	E638	L642	T646	L651	E652	Y653	E657	E658	Q665	L669	Y683	A684	LEU	ASP	GLY	GLN	ASN	ASP	ASN	LYS	SER	LYS	VAL	GLY	GLY	ILE	ALA	ARG	GLU	ILE	PHE	HIS	ASN	VAL	THR	ASN	PHE
------	------	------	------	------	------	------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.02Å 90.82Å 213.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.20 – 2.60 46.08 – 2.60	Depositor EDS
% Data completeness (in resolution range)	94.5 (31.20-2.60) 97.8 (46.08-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.08 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.205 , 0.271 0.211 , 0.275	Depositor DCC
R_{free} test set	2181 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	49.1	Xtriage
Anisotropy	0.397	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 48.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 43850 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7376	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.27% 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

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	C	0.43	0/2503	0.59	1/3376 (0.0%)
2	D	0.41	0/4998	0.55	0/6724
All	All	0.42	0/7501	0.57	1/10100 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	450	ARG	N-CA-C	-5.56	95.99	111.00

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	C	2464	0	2516	46	0
2	D	4909	0	4886	136	0
3	D	3	0	0	0	0
All	All	7376	0	7402	180	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 12.

All (180) 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:435:TYR:HB2	2:D:473:MSE:HE2	1.42	0.97
2:D:316:GLU:HG3	2:D:317:GLN:H	1.32	0.94
2:D:348:THR:HG22	2:D:349:GLN:HG3	1.57	0.87
2:D:546:THR:HG21	2:D:576:HIS:CE1	2.11	0.86
2:D:603:ALA:HB1	2:D:607:MSE:HG3	1.57	0.84
2:D:500:ASN:H	2:D:531:ASN:HD21	1.22	0.84
2:D:132:ASN:HD21	2:D:182:ILE:H	1.27	0.82
1:C:467:ALA:O	1:C:471:MSE:HG3	1.82	0.79
2:D:435:TYR:HB2	2:D:473:MSE:CE	2.13	0.78
2:D:435:TYR:CB	2:D:473:MSE:HE2	2.16	0.73
2:D:540:TYR:O	2:D:576:HIS:HD2	1.72	0.72
2:D:590:TYR:HD1	2:D:646:ILE:HD11	1.54	0.72
2:D:565:THR:O	2:D:569:GLU:HG3	1.89	0.72
2:D:622:MSE:SE	2:D:646:ILE:HD13	2.40	0.71
2:D:360:SER:HB2	2:D:409:ILE:HD11	1.73	0.71
2:D:117:LEU:HD11	2:D:123:TRP:HA	1.72	0.71
2:D:590:TYR:CD1	2:D:646:ILE:HD11	2.26	0.71
2:D:387:ILE:HG23	2:D:399:MSE:HE2	1.73	0.70
2:D:192:GLU:O	2:D:196:ILE:HG23	1.91	0.69
1:C:654:ASP:OD1	1:C:678:ARG:NH2	2.25	0.69
2:D:542:ASP:OD1	2:D:545:THR:HG23	1.94	0.68
2:D:101:TYR:CE2	2:D:104:LEU:HD22	2.28	0.68
2:D:191:SER:O	2:D:195:LYS:HG3	1.94	0.67
2:D:108:ILE:HG13	2:D:109:SER:N	2.10	0.67
1:C:439:THR:HG22	1:C:443:GLN:HE21	1.60	0.67
2:D:33:SER:O	2:D:37:GLU:HG3	1.95	0.66
2:D:562:GLU:OE2	2:D:564:GLN:HG3	1.96	0.66
2:D:132:ASN:ND2	2:D:182:ILE:H	1.92	0.66
2:D:30:GLY:HA3	2:D:31:SER:O	1.95	0.66
1:C:511:LEU:HD11	1:C:565:ILE:HG23	1.78	0.66
2:D:316:GLU:HG3	2:D:317:GLN:N	2.09	0.66
1:C:644:TYR:OH	1:C:676:GLU:OE2	2.13	0.65
2:D:315:GLU:HG2	2:D:316:GLU:N	2.12	0.65
2:D:431:ASN:O	2:D:473:MSE:HE1	1.97	0.65
2:D:117:LEU:HD11	2:D:123:TRP:CA	2.26	0.64
1:C:636:LEU:HD23	1:C:669:THR:HG22	1.79	0.63
1:C:639:ILE:HA	1:C:642:MSE:HE3	1.81	0.63
1:C:618:TYR:CZ	1:C:622:ARG:HD2	2.34	0.62
1:C:431:VAL:HG11	1:C:488:TYR:CD2	2.34	0.62
2:D:500:ASN:H	2:D:531:ASN:ND2	1.98	0.60
2:D:590:TYR:HD1	2:D:646:ILE:CD1	2.15	0.60
2:D:175:ARG:NH2	2:D:176:LEU:HD21	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:134:MSE:O	2:D:135:ILE:HD13	2.02	0.59
2:D:387:ILE:CG2	2:D:399:MSE:HE2	2.33	0.59
2:D:603:ALA:CB	2:D:607:MSE:HG3	2.33	0.58
1:C:450:ARG:O	1:C:452:LYS:N	2.35	0.58
2:D:590:TYR:CZ	2:D:642:LEU:HD22	2.38	0.58
2:D:514:LYS:HE2	2:D:516:SER:OG	2.03	0.58
2:D:546:THR:HG21	2:D:576:HIS:HE1	1.67	0.57
2:D:316:GLU:HG2	2:D:318:LYS:HD2	1.87	0.57
2:D:112:LEU:O	2:D:115:LYS:HE3	2.03	0.57
2:D:316:GLU:CG	2:D:317:GLN:H	2.10	0.57
2:D:53:LEU:HD11	2:D:56:ARG:NH2	2.20	0.56
2:D:105:GLN:HA	2:D:108:ILE:HG12	1.86	0.56
2:D:160:GLY:O	2:D:164:GLY:N	2.38	0.56
2:D:564:GLN:O	2:D:568:VAL:HG23	2.06	0.56
2:D:638:GLU:OE1	2:D:684:ALA:HB1	2.05	0.55
2:D:554:TYR:HA	2:D:559:ILE:HD12	1.87	0.55
2:D:108:ILE:HG13	2:D:109:SER:H	1.70	0.55
2:D:567:LYS:NZ	2:D:614:GLU:OE1	2.40	0.55
2:D:238:LYS:HG3	2:D:239:GLN:HE21	1.72	0.55
2:D:315:GLU:CG	2:D:316:GLU:H	2.18	0.54
1:C:425:ILE:HG13	1:C:426:GLN:N	2.22	0.54
2:D:101:TYR:HE2	2:D:104:LEU:HD22	1.73	0.54
2:D:683:TYR:O	2:D:684:ALA:HB3	2.08	0.54
2:D:315:GLU:HG2	2:D:316:GLU:H	1.72	0.54
2:D:529:ILE:HA	2:D:532:LEU:HD12	1.89	0.53
2:D:318:LYS:HG2	2:D:345:PHE:CE2	2.43	0.53
2:D:132:ASN:HD21	2:D:182:ILE:N	2.01	0.53
2:D:60:GLU:O	2:D:61:GLU:HB2	2.08	0.53
2:D:219:ALA:HB3	2:D:220:PRO:HD3	1.91	0.52
2:D:155:SER:OG	2:D:157:LYS:HB2	2.09	0.52
2:D:52:LEU:HD13	2:D:123:TRP:CH2	2.45	0.52
2:D:454:ARG:HG2	2:D:492:ASP:OD2	2.10	0.52
2:D:60:GLU:HG2	2:D:153:VAL:HG11	1.91	0.52
2:D:124:LEU:O	2:D:128:ILE:HG13	2.11	0.51
2:D:616:TRP:CZ2	2:D:658:GLU:HG3	2.46	0.51
2:D:459:MSE:CE	2:D:488:LEU:HD22	2.42	0.50
1:C:573:ILE:O	1:C:577:LEU:HB3	2.11	0.50
1:C:431:VAL:CG1	1:C:488:TYR:CD2	2.95	0.50
2:D:366:GLU:HG3	2:D:376:PHE:CE1	2.46	0.49
2:D:490:ILE:HD13	2:D:553:ILE:HG12	1.95	0.49
2:D:360:SER:HB3	2:D:405:SER:OG	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:622:MSE:SE	2:D:646:ILE:CD1	3.11	0.49
1:C:471:MSE:HG2	2:D:651:LEU:CB	2.43	0.48
2:D:359:PHE:HE2	2:D:384:VAL:CG1	2.26	0.48
1:C:511:LEU:HD11	1:C:565:ILE:CG2	2.43	0.48
2:D:112:LEU:O	2:D:115:LYS:HB2	2.13	0.48
1:C:497:LEU:HD22	1:C:500:MSE:HE2	1.96	0.48
2:D:631:VAL:O	2:D:632:ASP:HB2	2.12	0.48
1:C:441:PHE:CE1	1:C:461:PHE:CD1	3.02	0.48
1:C:366:GLN:OE1	1:C:366:GLN:HA	2.14	0.48
2:D:579:TYR:HA	2:D:582:VAL:HB	1.95	0.48
1:C:523:ASN:OD1	1:C:581:ASN:ND2	2.47	0.48
2:D:30:GLY:CA	2:D:31:SER:HB2	2.44	0.47
2:D:143:SER:HA	2:D:146:TRP:CE2	2.50	0.47
2:D:345:PHE:O	2:D:348:THR:HB	2.14	0.47
2:D:459:MSE:HE1	2:D:488:LEU:HD22	1.97	0.47
2:D:562:GLU:HG2	2:D:562:GLU:O	2.14	0.47
2:D:658:GLU:HA	2:D:658:GLU:OE1	2.15	0.47
2:D:665:GLN:HE21	2:D:669:LEU:HD11	1.80	0.47
2:D:514:LYS:H	2:D:517:ASN:ND2	2.12	0.47
2:D:612:GLY:HA3	2:D:653:VAL:O	2.14	0.47
2:D:522:ARG:HG2	2:D:522:ARG:O	2.16	0.46
1:C:613:ASN:O	1:C:619:ARG:NH1	2.49	0.46
2:D:168:PRO:HG3	2:D:220:PRO:HB2	1.97	0.46
2:D:134:MSE:C	2:D:135:ILE:HD13	2.36	0.46
2:D:315:GLU:CG	2:D:316:GLU:N	2.73	0.46
2:D:126:GLU:O	2:D:130:ILE:HG13	2.14	0.46
2:D:101:TYR:HD2	2:D:101:TYR:O	1.99	0.46
2:D:350:GLU:OE1	2:D:358:HIS:HD2	1.99	0.46
2:D:61:GLU:O	2:D:62:VAL:C	2.54	0.46
1:C:483:TYR:CD1	1:C:503:ASN:ND2	2.84	0.46
2:D:117:LEU:HD11	2:D:123:TRP:CB	2.46	0.46
1:C:358:GLU:HG2	1:C:429:VAL:HG12	1.98	0.45
1:C:582:ILE:HD13	1:C:596:LEU:HD22	1.97	0.45
2:D:57:GLU:C	2:D:59:GLU:H	2.20	0.45
1:C:558:TYR:HD1	1:C:566:LEU:HD11	1.82	0.45
1:C:562:ASP:OD1	1:C:564:THR:HB	2.16	0.45
1:C:468:PHE:CE2	1:C:472:VAL:HG21	2.52	0.45
2:D:106:GLU:O	2:D:110:LYS:HB2	2.16	0.45
1:C:361:GLU:O	1:C:429:VAL:HG23	2.17	0.45
1:C:579:ILE:HD13	1:C:624:LYS:HD3	1.97	0.44
2:D:31:SER:CB	2:D:34:LYS:HG2	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:351:GLU:HG2	2:D:352:GLU:N	2.32	0.44
1:C:627:ILE:HD13	1:C:647:GLU:O	2.16	0.44
2:D:407:LEU:CD1	2:D:434:ILE:HD12	2.47	0.44
2:D:435:TYR:CA	2:D:473:MSE:HE2	2.47	0.44
1:C:431:VAL:CG1	1:C:488:TYR:HD2	2.30	0.44
2:D:52:LEU:HD13	2:D:123:TRP:HH2	1.80	0.44
1:C:558:TYR:CD1	1:C:566:LEU:HD11	2.53	0.44
2:D:490:ILE:HG13	2:D:491:SER:N	2.31	0.44
1:C:343:MSE:HE1	1:C:441:PHE:CD2	2.53	0.44
2:D:536:ASN:HA	2:D:537:PRO:HD3	1.80	0.44
2:D:543:LEU:N	2:D:544:PRO:CD	2.80	0.44
1:C:683:GLU:O	1:C:684:ALA:C	2.56	0.44
1:C:524:ASP:HB3	1:C:540:ILE:HD13	2.00	0.44
2:D:488:LEU:HD23	2:D:488:LEU:HA	1.75	0.44
1:C:625:LEU:HD23	1:C:625:LEU:HA	1.83	0.43
1:C:640:LEU:HG	1:C:644:TYR:CE2	2.53	0.43
1:C:471:MSE:HG2	2:D:651:LEU:HB3	1.99	0.43
2:D:366:GLU:HG3	2:D:376:PHE:CZ	2.54	0.43
1:C:518:ILE:HD13	1:C:573:ILE:HD11	1.99	0.43
1:C:347:ARG:HA	1:C:347:ARG:NE	2.34	0.43
2:D:117:LEU:HD11	2:D:123:TRP:HB2	2.01	0.43
1:C:636:LEU:HD12	1:C:636:LEU:HA	1.79	0.43
1:C:636:LEU:HD21	1:C:670:ARG:HA	2.01	0.43
2:D:263:VAL:O	2:D:264:LYS:HG3	2.19	0.42
2:D:192:GLU:HA	2:D:195:LYS:HB2	2.01	0.42
1:C:453:VAL:HG12	1:C:457:TYR:HB2	2.00	0.42
2:D:101:TYR:O	2:D:104:LEU:HB2	2.19	0.42
1:C:340:PHE:HE1	1:C:453:VAL:HG13	1.85	0.42
1:C:519:TYR:CZ	1:C:523:ASN:ND2	2.88	0.42
2:D:249:TYR:O	2:D:252:PHE:HB3	2.20	0.42
2:D:598:GLU:O	2:D:599:ILE:C	2.57	0.42
2:D:53:LEU:HD11	2:D:56:ARG:HH21	1.84	0.42
2:D:143:SER:O	2:D:147:GLU:HB2	2.20	0.42
2:D:255:LEU:HD23	2:D:255:LEU:HA	1.81	0.42
1:C:640:LEU:HA	1:C:640:LEU:HD12	1.82	0.41
2:D:46:LEU:CD1	2:D:135:ILE:HD11	2.50	0.41
2:D:39:LEU:HA	2:D:39:LEU:HD12	1.82	0.41
2:D:213:VAL:HG12	2:D:214:ILE:HD13	2.02	0.41
2:D:549:LEU:O	2:D:553:ILE:HG13	2.20	0.41
2:D:456:GLU:HA	2:D:457:PRO:HD3	1.94	0.41
2:D:218:LEU:HD11	2:D:231:PHE:CE1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:261:LYS:HE2	2:D:261:LYS:HB3	1.78	0.41
2:D:30:GLY:HA2	2:D:31:SER:HB2	2.03	0.41
2:D:574:VAL:HG13	2:D:618:THR:HG21	2.01	0.41
2:D:683:TYR:O	2:D:684:ALA:CB	2.68	0.41
2:D:140:LEU:O	2:D:141:LEU:C	2.58	0.41
2:D:213:VAL:O	2:D:217:GLU:HB2	2.20	0.41
2:D:55:LEU:HD23	2:D:55:LEU:HA	1.80	0.41
2:D:411:LEU:HG	2:D:419:ILE:HD13	2.03	0.41
1:C:341:GLU:O	1:C:345:GLN:HG3	2.21	0.41
2:D:411:LEU:HG	2:D:419:ILE:CD1	2.51	0.40
2:D:562:GLU:C	2:D:564:GLN:H	2.24	0.40
2:D:562:GLU:C	2:D:564:GLN:N	2.72	0.40
2:D:562:GLU:CD	2:D:562:GLU:H	2.24	0.40
2:D:311:GLU:O	2:D:312:ILE:HD13	2.21	0.40
2:D:522:ARG:NH1	2:D:557:LEU:O	2.55	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	C	291/357 (82%)	278 (96%)	11 (4%)	2 (1%)	26	51
2	D	584/709 (82%)	548 (94%)	31 (5%)	5 (1%)	21	42
All	All	875/1066 (82%)	826 (94%)	42 (5%)	7 (1%)	24	46

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	448	ILE
1	C	450	ARG
2	D	316	GLU

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Mol	Chain	Res	Type
2	D	31	SER
2	D	61	GLU
2	D	151	SER
2	D	603	ALA

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	C	281/330 (85%)	274 (98%)	7 (2%)	55	81
2	D	555/641 (87%)	526 (95%)	29 (5%)	29	54
All	All	836/971 (86%)	800 (96%)	36 (4%)	35	64

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	343	MSE
1	C	363	VAL
1	C	448	ILE
1	C	458	LEU
1	C	501	GLU
1	C	570	ILE
1	C	650	LEU
2	D	38	ILE
2	D	46	LEU
2	D	58	LEU
2	D	101	TYR
2	D	117	LEU
2	D	122	GLU
2	D	140	LEU
2	D	141	LEU
2	D	147	GLU
2	D	148	THR
2	D	183	ASN
2	D	193	LEU

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Mol	Chain	Res	Type
2	D	196	ILE
2	D	223	SER
2	D	246	ASP
2	D	264	LYS
2	D	268	GLU
2	D	303	SER
2	D	311	GLU
2	D	323	LEU
2	D	341	PHE
2	D	348	THR
2	D	427	ASP
2	D	456	GLU
2	D	524	CYS
2	D	565	THR
2	D	582	VAL
2	D	622	MSE
2	D	657	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	443	GLN
1	C	563	ASN
2	D	132	ASN
2	D	239	GLN
2	D	358	HIS
2	D	430	ASN
2	D	517	ASN
2	D	531	ASN
2	D	576	HIS
2	D	665	GLN

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	C	289/357 (80%)	-0.02	8 (2%) 56 49	30, 51, 106, 149	0
2	D	580/709 (81%)	0.01	15 (2%) 59 53	35, 59, 102, 145	0
All	All	869/1066 (81%)	-0.00	23 (2%) 59 53	30, 56, 103, 149	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	102	HIS	5.5
2	D	38	ILE	4.0
1	C	363	VAL	3.4
1	C	425	ILE	3.3
2	D	36	LEU	2.9
1	C	453	VAL	2.8
2	D	30	GLY	2.7
2	D	39	LEU	2.6
1	C	365	VAL	2.6
2	D	101	TYR	2.5
2	D	31	SER	2.5
2	D	104	LEU	2.5
1	C	335	TYR	2.5
1	C	424	PRO	2.5
1	C	364	SER	2.4
2	D	58	LEU	2.4
2	D	52	LEU	2.3
2	D	42	LEU	2.2
2	D	62	VAL	2.2
2	D	105	GLN	2.2
1	C	362	LEU	2.1
2	D	55	LEU	2.1
2	D	40	CYS	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.