Full wwPDB X-ray Structure Validation Report (i)

Sep 25, 2023 – 06:00 AM EDT

PDB ID	:	5VNK
Title	:	Crystal structure of Sec23a/Sec24a/Sec22 complexed with a C-terminal LL sorting motif
Authors	:	Ma, W.; Goldberg, J.
Deposited on	:	2017-04-30
Resolution	:	2.55 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

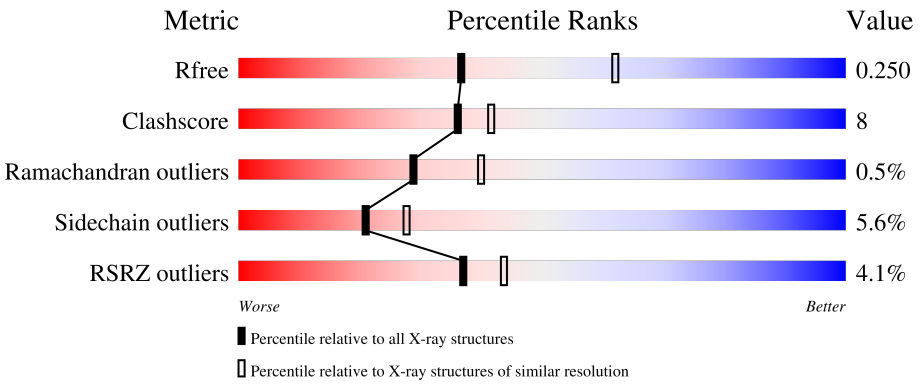
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.55 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}	130704	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	764	4%, 74%, 16%, •, 7%
2	B	748	2%, 78%, 17%, •, •
3	C	157	8%, 62%, 22%, •, •, 14%
4	D	5	60%, 20%, 40%, 40%

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
5	ZN	A	801	-	-	X	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 12544 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 Protein transport protein Sec23A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	708	5620	3582	967	1031	40	0	0	0

- Molecule 2 is a protein called Protein transport protein Sec24A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	729	5754	3669	980	1071	34	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1056	ALA	ARG	conflict	UNP O95486

- Molecule 3 is a protein called Vesicle-trafficking protein SEC22b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	135	1087	699	177	203	8	0	0	0

- Molecule 4 is a protein called C-terminal LL.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	D	3	22	15	3	4	0	0	0

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
5	A	1	1	1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
5	B	1	1	1	0	0

- Molecule 6 is water.

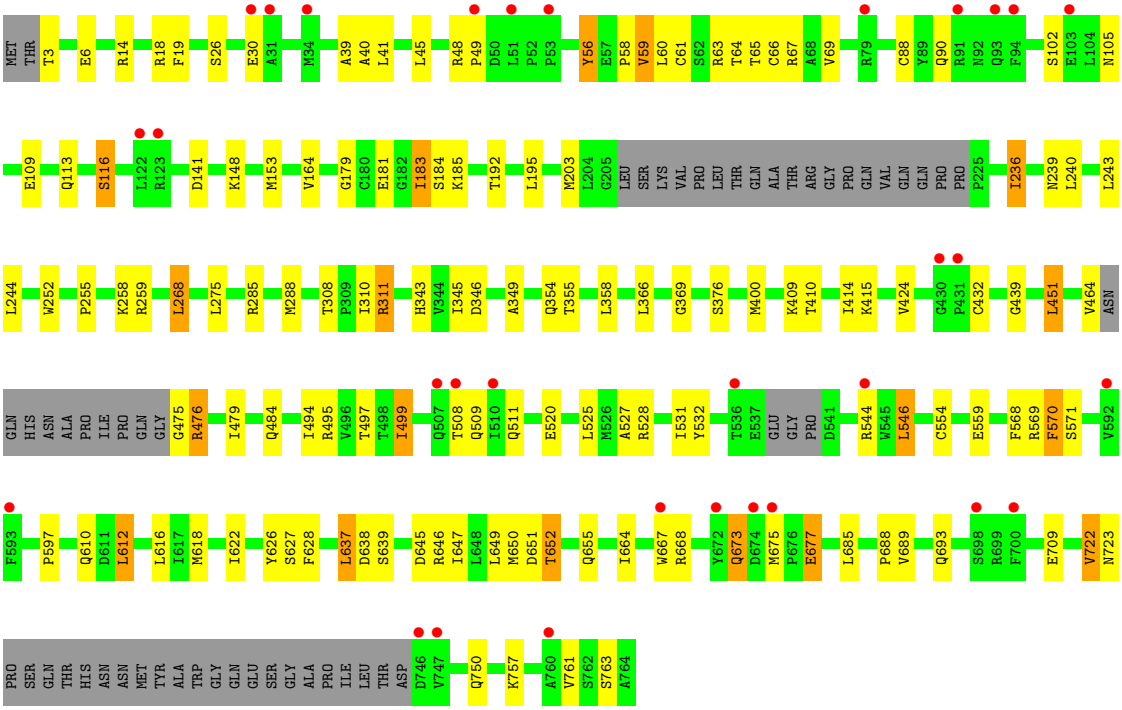
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	21	21	21	0	0
			Total	O		
6	B	35	35	35	0	0
			Total	O		
6	C	3	3	3	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

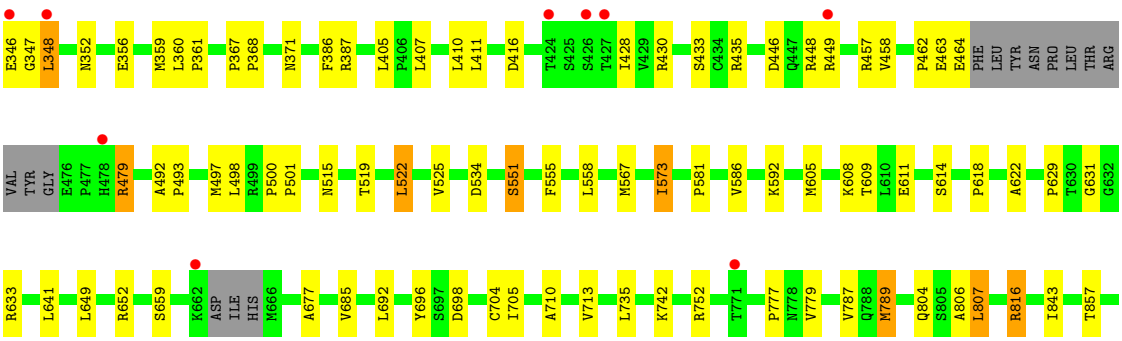
• Molecule 1: Protein transport protein Sec23A

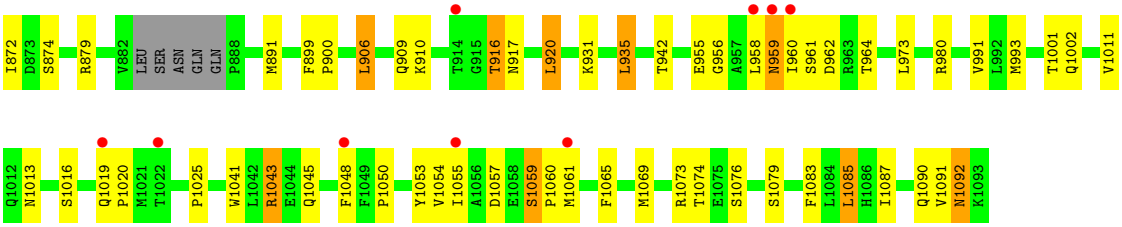
Chain A: 4% 74% 16% 7%



• Molecule 2: Protein transport protein Sec24A

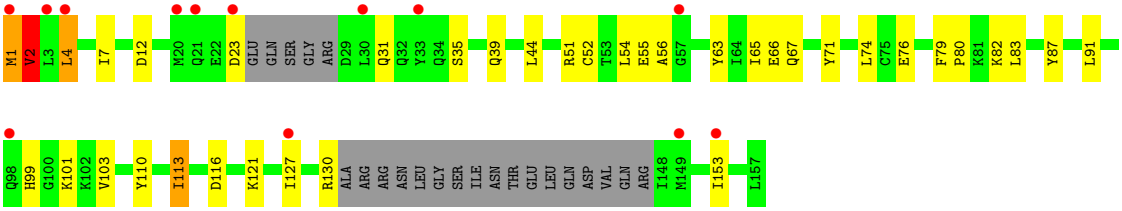
Chain B: 2% 78% 17%





- Molecule 3: Vesicle-trafficking protein SEC22b

Chain C:



- Molecule 4: C-terminal LL

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	148.18Å 97.61Å 130.55Å 90.00° 90.03° 90.00°	Depositor
Resolution (Å)	130.55 – 2.55 130.55 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.1 (130.55-2.55) 99.1 (130.55-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.74 (at 2.55Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.196 , 0.251 0.197 , 0.250	Depositor DCC
R_{free} test set	3041 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	57.3	Xtriage
Anisotropy	0.128	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 48.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.018 for -h,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	12544	wwPDB-VP
Average B, all atoms (Å^2)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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: ZN

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	0.42	0/5751	0.51	0/7786
2	B	0.34	0/5877	0.50	0/7989
3	C	0.29	0/1106	0.50	0/1489
4	D	0.30	0/21	0.89	0/26
All	All	0.38	0/12755	0.51	0/17290

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	5620	0	5567	89	0
2	B	5754	0	5795	88	0
3	C	1087	0	1091	27	0
4	D	22	0	23	2	0
5	A	1	0	0	2	0
5	B	1	0	0	0	0
6	A	21	0	0	2	0
6	B	35	0	0	10	0
6	C	3	0	0	0	0
All	All	12544	0	12476	200	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 8.

All (200) 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:61:CYS:SG	5:A:801:ZN:ZN	1.52	0.98
2:B:551:SER:OG	2:B:611:GLU:OE1	1.94	0.85
3:C:1:MET:H1	3:C:76:GLU:HB2	1.39	0.83
3:C:1:MET:N	3:C:1:MET:HE3	1.95	0.82
1:A:60:LEU:HD21	1:A:69:VAL:HG22	1.62	0.80
2:B:1057:ASP:HB2	2:B:1060:PRO:HD3	1.64	0.79
1:A:60:LEU:HD23	1:A:69:VAL:HA	1.65	0.78
1:A:179:GLY:HA2	1:A:239:ASN:HD22	1.48	0.78
1:A:255:PRO:HG2	1:A:258:LYS:HG3	1.65	0.77
2:B:629:PRO:HG3	3:C:23:ASP:HB2	1.66	0.76
2:B:961:SER:OG	2:B:962:ASP:OD1	2.04	0.75
1:A:259:ARG:NH2	1:A:308:THR:O	2.19	0.75
2:B:779:VAL:HG21	2:B:807:LEU:HD21	1.72	0.72
1:A:61:CYS:HG	5:A:801:ZN:ZN	0.40	0.71
1:A:63:ARG:NH1	1:A:90:GLN:HB2	2.07	0.70
1:A:432:CYS:SG	6:A:916:HOH:O	2.50	0.69
3:C:1:MET:N	3:C:76:GLU:HB2	2.07	0.69
2:B:1074:THR:HG23	2:B:1076:SER:H	1.58	0.69
1:A:647:ILE:HD11	1:A:664:ILE:HG21	1.75	0.68
1:A:647:ILE:HG21	1:A:688:PRO:HG3	1.77	0.67
2:B:1019:GLN:HB3	2:B:1020:PRO:HD3	1.77	0.66
1:A:60:LEU:CD2	1:A:69:VAL:HG22	2.25	0.66
2:B:804:GLN:OE1	6:B:1201:HOH:O	2.14	0.65
1:A:3:THR:HG22	1:A:6:GLU:H	1.62	0.65
2:B:1025:PRO:O	6:B:1202:HOH:O	2.14	0.65
2:B:958:LEU:HD12	2:B:958:LEU:N	2.12	0.64
3:C:1:MET:H1	3:C:1:MET:HE3	1.62	0.64
2:B:806:ALA:HB3	4:D:358:LEU:HD11	1.80	0.63
3:C:4:LEU:HB3	3:C:74:LEU:HB3	1.81	0.63
1:A:48:ARG:HB2	1:A:49:PRO:HD2	1.81	0.63
1:A:56:TYR:OH	1:A:109:GLU:OE2	2.17	0.63
1:A:508:THR:HG22	1:A:511:GLN:OE1	1.99	0.62
2:B:1054:VAL:O	2:B:1054:VAL:HG12	1.99	0.62
1:A:102:SER:O	1:A:105:ASN:O	2.17	0.62
3:C:110:TYR:HB3	3:C:113:ILE:HG23	1.82	0.62
2:B:492:ALA:HB3	2:B:816:ARG:HB3	1.82	0.62
2:B:371:ASN:OD1	6:B:1204:HOH:O	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:627:SER:HB3	1:A:646:ARG:HG3	1.82	0.61
2:B:891:MET:O	6:B:1203:HOH:O	2.15	0.60
1:A:626:TYR:HB2	1:A:647:ILE:HB	1.82	0.60
1:A:61:CYS:O	1:A:67:ARG:HA	2.01	0.60
2:B:430:ARG:HD3	2:B:435:ARG:HH21	1.66	0.60
1:A:610:GLN:HG3	1:A:618:MET:HE1	1.82	0.59
1:A:183:ILE:HD13	2:B:605:MET:CE	2.32	0.59
2:B:959:ASN:O	2:B:959:ASN:ND2	2.35	0.59
2:B:962:ASP:OD1	2:B:962:ASP:N	2.27	0.59
2:B:1055:ILE:O	2:B:1057:ASP:N	2.35	0.59
4:D:357:LEU:HD23	4:D:357:LEU:H	1.67	0.59
1:A:475:GLY:O	1:A:476:ARG:HB2	2.03	0.58
2:B:348:LEU:HD12	2:B:348:LEU:O	2.04	0.57
1:A:61:CYS:SG	1:A:88:CYS:SG	3.02	0.57
3:C:113:ILE:O	3:C:116:ASP:HB2	2.04	0.57
1:A:26:SER:H	1:A:509:GLN:NE2	2.04	0.56
1:A:673:GLN:HG2	1:A:685:LEU:HD12	1.85	0.56
3:C:35:SER:O	3:C:39:GLN:HG2	2.06	0.56
2:B:872:ILE:HD12	2:B:1090:GLN:HB2	1.88	0.56
2:B:931:LYS:NZ	6:B:1207:HOH:O	2.31	0.56
2:B:1043:ARG:HB3	2:B:1050:PRO:HD2	1.88	0.55
1:A:67:ARG:O	1:A:409:LYS:NZ	2.38	0.55
2:B:405:LEU:HD22	2:B:843:ILE:HD13	1.89	0.55
2:B:1073:ARG:NH2	6:B:1209:HOH:O	2.40	0.55
1:A:410:THR:HB	1:A:414:ILE:HB	1.88	0.54
2:B:955:GLU:N	2:B:955:GLU:OE1	2.40	0.54
1:A:354:GLN:HE22	1:A:597:PRO:HD2	1.72	0.54
1:A:116:SER:OG	1:A:497:THR:OG1	1.81	0.54
3:C:54:LEU:HD13	3:C:153:ILE:HD13	1.89	0.54
2:B:906:LEU:HD13	2:B:942:THR:HG21	1.90	0.54
1:A:195:LEU:HD22	1:A:203:MET:HE1	1.90	0.54
1:A:45:LEU:HD11	1:A:451:LEU:HD13	1.89	0.53
2:B:1083:PHE:O	2:B:1087:ILE:HG12	2.08	0.53
3:C:7:ILE:HD12	3:C:71:TYR:CD2	2.43	0.53
2:B:909:GLN:HG2	2:B:910:LYS:N	2.24	0.53
1:A:439:GLY:HA2	1:A:532:TYR:CZ	2.43	0.53
1:A:415:LYS:HD2	1:A:464:VAL:HG21	1.89	0.53
1:A:637:LEU:HD22	1:A:722:VAL:HA	1.92	0.52
1:A:56:TYR:HH	1:A:109:GLU:CD	2.12	0.52
2:B:387:ARG:HD2	2:B:935:LEU:HD12	1.92	0.52
2:B:416:ASP:OD1	2:B:742:LYS:NZ	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:558:LEU:HB2	2:B:586:VAL:HG11	1.90	0.52
2:B:872:ILE:HD13	2:B:1087:ILE:HG23	1.92	0.52
3:C:80:PRO:HB3	3:C:82:LYS:HE2	1.93	0.51
3:C:44:LEU:HD13	3:C:65:ILE:HD11	1.93	0.51
1:A:479:ILE:HB	1:A:499:ILE:HD11	1.92	0.50
2:B:958:LEU:N	2:B:958:LEU:CD1	2.74	0.50
1:A:366:LEU:HD22	1:A:424:VAL:HG22	1.93	0.50
2:B:980:ARG:NH2	6:B:1211:HOH:O	2.43	0.50
2:B:961:SER:OG	2:B:962:ASP:N	2.45	0.50
2:B:446:ASP:C	2:B:448:ARG:H	2.14	0.50
1:A:58:PRO:O	1:A:59:VAL:HG23	2.11	0.49
2:B:410:LEU:HD22	2:B:935:LEU:HG	1.94	0.49
3:C:2:VAL:O	3:C:127:ILE:HD11	2.12	0.49
2:B:652:ARG:HB2	2:B:696:TYR:CE2	2.48	0.48
1:A:311:ARG:HD3	1:A:311:ARG:H	1.79	0.48
1:A:39:ALA:HB3	1:A:525:LEU:HD13	1.95	0.48
1:A:63:ARG:O	1:A:64:THR:C	2.50	0.48
1:A:58:PRO:O	1:A:59:VAL:CB	2.62	0.48
2:B:752:ARG:NH1	6:B:1201:HOH:O	2.28	0.48
1:A:439:GLY:HA2	1:A:532:TYR:CE2	2.49	0.47
1:A:652:THR:HG22	1:A:655:GLN:H	1.80	0.47
2:B:1073:ARG:HB3	2:B:1079:SER:HB3	1.95	0.47
2:B:649:LEU:HD13	2:B:698:ASP:HB3	1.97	0.47
1:A:195:LEU:HD22	1:A:203:MET:CE	2.44	0.47
1:A:622:ILE:HG13	1:A:651:ASP:HB3	1.97	0.47
1:A:14:ARG:O	1:A:48:ARG:NH1	2.48	0.47
1:A:554:CYS:HA	1:A:570:PHE:CZ	2.49	0.47
2:B:433:SER:OG	2:B:457:ARG:HD3	2.15	0.47
2:B:631:GLY:HA2	2:B:685:VAL:HG22	1.97	0.47
3:C:56:ALA:HB2	3:C:153:ILE:HG21	1.95	0.47
1:A:757:LYS:O	1:A:761:VAL:HG22	2.15	0.47
1:A:628:PHE:HE2	1:A:667:TRP:HZ3	1.63	0.46
1:A:275:LEU:HB3	1:A:343:HIS:CE1	2.50	0.46
2:B:462:PRO:O	2:B:464:GLU:N	2.48	0.46
2:B:573:ILE:HG23	2:B:618:PRO:HG2	1.96	0.46
1:A:184:SER:O	1:A:184:SER:OG	2.31	0.46
2:B:677:ALA:HB2	2:B:705:ILE:HA	1.98	0.46
1:A:185:LYS:HB3	2:B:567:MET:HB3	1.98	0.46
1:A:476:ARG:HG2	1:A:476:ARG:HH11	1.81	0.46
2:B:899:PHE:HB3	2:B:900:PRO:HD3	1.97	0.46
3:C:52:CYS:HB3	3:C:63:TYR:CE2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ARG:HH12	1:A:90:GLN:HB2	1.80	0.46
2:B:879:ARG:CZ	2:B:1092:ASN:HB3	2.46	0.46
2:B:1019:GLN:CB	2:B:1020:PRO:HD3	2.46	0.46
1:A:349:ALA:HB1	1:A:355:THR:HG21	1.98	0.45
3:C:7:ILE:HD12	3:C:71:TYR:HD2	1.80	0.45
1:A:638:ASP:OD1	1:A:639:SER:N	2.50	0.45
1:A:559:GLU:O	1:A:568:PHE:HA	2.16	0.45
2:B:534:ASP:OD1	2:B:592:LYS:NZ	2.32	0.45
1:A:236:ILE:HD13	1:A:236:ILE:HA	1.73	0.45
2:B:609:THR:HG22	2:B:611:GLU:H	1.80	0.45
3:C:1:MET:N	3:C:1:MET:CE	2.72	0.45
2:B:973:LEU:HA	2:B:1069:MET:HE1	1.99	0.45
2:B:479:ARG:HD2	2:B:479:ARG:HA	1.74	0.45
1:A:527:ALA:O	1:A:531:ILE:HG12	2.17	0.44
2:B:916:THR:HG22	2:B:917:ASN:H	1.82	0.44
2:B:960:ILE:O	2:B:960:ILE:HG22	2.17	0.44
2:B:956:GLY:C	2:B:958:LEU:HD12	2.38	0.44
3:C:1:MET:HE3	3:C:1:MET:CA	2.39	0.44
1:A:345:ILE:O	1:A:369:GLY:HA3	2.17	0.44
3:C:55:GLU:O	3:C:153:ILE:HG22	2.18	0.44
1:A:19:PHE:CE1	1:A:40:ALA:HB2	2.52	0.44
3:C:1:MET:HA	3:C:79:PHE:HB2	2.00	0.44
1:A:48:ARG:NH2	6:A:906:HOH:O	2.50	0.43
1:A:63:ARG:O	1:A:66:CYS:N	2.45	0.43
2:B:633:ARG:NH2	6:B:1215:HOH:O	2.50	0.43
2:B:457:ARG:HG3	2:B:458:VAL:N	2.32	0.43
2:B:525:VAL:HG22	2:B:735:LEU:HD11	1.99	0.43
2:B:1041:TRP:O	2:B:1045:GLN:HG2	2.18	0.43
3:C:1:MET:N	3:C:76:GLU:OE1	2.51	0.43
1:A:113:GLN:H	1:A:113:GLN:HG3	1.50	0.43
2:B:407:LEU:HG	2:B:789:MET:HG3	2.00	0.43
2:B:704:CYS:SG	6:B:1234:HOH:O	2.62	0.43
3:C:80:PRO:HB2	3:C:83:LEU:HG	2.00	0.43
2:B:346:GLU:HB3	2:B:347:GLY:H	1.53	0.43
1:A:722:VAL:HG13	1:A:723:ASN:N	2.34	0.43
2:B:1085:LEU:HD12	2:B:1085:LEU:HA	1.86	0.43
1:A:675:MET:HB3	1:A:677:GLU:HG3	2.00	0.43
1:A:310:ILE:HG22	1:A:311:ARG:HD3	2.00	0.42
2:B:1060:PRO:O	2:B:1061:MET:HB2	2.19	0.42
1:A:546:LEU:HD12	1:A:546:LEU:HA	1.87	0.42
1:A:484:GLN:HG2	1:A:494:ILE:HG13	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:689:VAL:O	1:A:693:GLN:HG2	2.20	0.42
3:C:12:ASP:OD1	3:C:12:ASP:N	2.44	0.42
3:C:99:HIS:O	3:C:103:VAL:HG23	2.19	0.42
1:A:58:PRO:O	1:A:59:VAL:HB	2.19	0.42
2:B:1011:VAL:HG12	2:B:1013:ASN:H	1.84	0.42
2:B:920:LEU:HD12	2:B:920:LEU:HA	1.90	0.42
3:C:87:TYR:CZ	3:C:91:LEU:HD11	2.54	0.42
1:A:252:TRP:CE2	2:B:581:PRO:HD3	2.55	0.42
1:A:60:LEU:CD2	1:A:69:VAL:HA	2.43	0.42
1:A:268:LEU:HG	1:A:288:MET:SD	2.59	0.42
3:C:51:ARG:NH2	3:C:66:GLU:OE2	2.53	0.42
1:A:285:ARG:NE	1:A:346:ASP:OD2	2.45	0.41
2:B:710:ALA:HB3	2:B:777:PRO:HD2	2.01	0.41
1:A:358:LEU:HD22	1:A:597:PRO:HB3	2.02	0.41
1:A:63:ARG:O	1:A:65:THR:N	2.53	0.41
2:B:522:LEU:HD23	2:B:522:LEU:HA	1.90	0.41
2:B:428:ILE:HD13	2:B:493:PRO:HD3	2.03	0.41
2:B:360:LEU:HD12	2:B:361:PRO:HD2	2.03	0.41
2:B:555:PHE:HZ	2:B:622:ALA:HB1	1.85	0.41
1:A:185:LYS:CB	2:B:567:MET:HB3	2.51	0.41
1:A:56:TYR:OH	1:A:109:GLU:CD	2.59	0.41
1:A:649:LEU:HD12	1:A:650:MET:N	2.36	0.41
2:B:515:ASN:O	2:B:519:THR:HG23	2.21	0.41
2:B:991:VAL:HG21	2:B:1053:TYR:HE1	1.85	0.41
1:A:520:GLU:HB3	1:A:616:LEU:HD11	2.03	0.41
1:A:645:ASP:HA	1:A:664:ILE:HG12	2.03	0.41
2:B:367:PRO:HA	2:B:368:PRO:HD3	1.96	0.41
2:B:910:LYS:HE2	2:B:910:LYS:HB2	1.89	0.41
2:B:956:GLY:O	2:B:958:LEU:CD1	2.69	0.41
1:A:354:GLN:NE2	1:A:597:PRO:HD2	2.35	0.40
2:B:1059:SER:N	2:B:1060:PRO:CD	2.83	0.40
2:B:1087:ILE:O	2:B:1091:VAL:HG23	2.21	0.40
2:B:348:LEU:O	2:B:348:LEU:CG	2.70	0.40
2:B:446:ASP:HB2	2:B:449:ARG:HB2	2.03	0.40
1:A:148:LYS:HE3	1:A:244:LEU:O	2.21	0.40
1:A:18:ARG:NH1	1:A:612:LEU:HD22	2.36	0.40
1:A:685:LEU:HD23	1:A:685:LEU:HA	1.88	0.40
1:A:750:GLN:OE1	1:A:750:GLN:N	2.52	0.40
2:B:352:ASN:O	2:B:356:GLU:HG2	2.20	0.40
2:B:500:PRO:HA	2:B:501:PRO:HD3	2.00	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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	698/764 (91%)	667 (96%)	28 (4%)	3 (0%)	34	46
2	B	721/748 (96%)	680 (94%)	38 (5%)	3 (0%)	34	46
3	C	129/157 (82%)	118 (92%)	9 (7%)	2 (2%)	9	12
4	D	1/5 (20%)	0	1 (100%)	0	100	100
All	All	1549/1674 (92%)	1465 (95%)	76 (5%)	8 (0%)	29	40

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	4	LEU
1	A	59	VAL
2	B	463	GLU
1	A	476	ARG
2	B	1065	PHE
3	C	2	VAL
1	A	181	GLU
2	B	1059	SER

5.3.2 Protein sidechains (i)

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	617/666 (93%)	583 (94%)	34 (6%)	21	29
2	B	658/678 (97%)	622 (94%)	36 (6%)	21	29

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	C	119/138 (86%)	111 (93%)	8 (7%)	16 21
4	D	2/5 (40%)	2 (100%)	0	100 100
All	All	1396/1487 (94%)	1318 (94%)	78 (6%)	21 28

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLU
1	A	41	LEU
1	A	56	TYR
1	A	116	SER
1	A	141	ASP
1	A	153	MET
1	A	164	VAL
1	A	183	ILE
1	A	192	THR
1	A	236	ILE
1	A	240	LEU
1	A	243	LEU
1	A	268	LEU
1	A	311	ARG
1	A	376	SER
1	A	400	MET
1	A	451	LEU
1	A	495	ARG
1	A	499	ILE
1	A	528	ARG
1	A	544	ARG
1	A	546	LEU
1	A	569	ARG
1	A	570	PHE
1	A	571	SER
1	A	612	LEU
1	A	637	LEU
1	A	652	THR
1	A	668	ARG
1	A	673	GLN
1	A	677	GLU
1	A	709	GLU
1	A	722	VAL
1	A	763	SER

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Mol	Chain	Res	Type
2	B	348	LEU
2	B	359	MET
2	B	386	PHE
2	B	411	LEU
2	B	479	ARG
2	B	497	MET
2	B	498	LEU
2	B	522	LEU
2	B	551	SER
2	B	573	ILE
2	B	608	LYS
2	B	614	SER
2	B	641	LEU
2	B	659	SER
2	B	692	LEU
2	B	713	VAL
2	B	787	VAL
2	B	789	MET
2	B	807	LEU
2	B	816	ARG
2	B	857	THR
2	B	874	SER
2	B	906	LEU
2	B	916	THR
2	B	920	LEU
2	B	935	LEU
2	B	959	ASN
2	B	964	THR
2	B	993	MET
2	B	1001	THR
2	B	1002	GLN
2	B	1016	SER
2	B	1043	ARG
2	B	1048	PHE
2	B	1085	LEU
2	B	1092	ASN
3	C	1	MET
3	C	2	VAL
3	C	31	GLN
3	C	67	GLN
3	C	101	LYS
3	C	113	ILE

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Mol	Chain	Res	Type
3	C	121	LYS
3	C	130	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	239	ASN
1	A	256	GLN
2	B	729	GLN
2	B	788	GLN
2	B	959	ASN
2	B	1019	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	708/764 (92%)	0.41	31 (4%)	34	41	38, 69, 112, 158	0
2	B	729/748 (97%)	0.30	18 (2%)	57	63	32, 58, 102, 140	0
3	C	135/157 (85%)	0.70	13 (9%)	8	10	53, 88, 136, 155	0
4	D	3/5 (60%)	5.29	3 (100%)	0	0	140, 140, 141, 143	0
All	All	1575/1674 (94%)	0.39	65 (4%)	37	44	32, 65, 114, 158	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1055	ILE	7.5
2	B	959	ASN	7.2
3	C	1	MET	6.7
4	D	358	LEU	6.5
3	C	153	ILE	6.0
4	D	357	LEU	5.0
4	D	356	SER	4.4
3	C	23	ASP	4.3
1	A	747	VAL	4.3
2	B	1061	MET	4.2
1	A	49	PRO	3.9
1	A	31	ALA	3.8
2	B	426	SER	3.7
3	C	127	ILE	3.6
2	B	1019	GLN	3.6
1	A	93	GLN	3.6
1	A	746	ASP	3.4
3	C	57	GLY	3.4
2	B	958	LEU	3.4
1	A	30	GLU	3.4
1	A	698	SER	3.4

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Mol	Chain	Res	Type	RSRZ
3	C	3	LEU	3.2
3	C	4	LEU	3.2
1	A	123	ARG	3.2
2	B	960	ILE	3.1
1	A	430	GLY	3.0
1	A	34	MET	2.8
2	B	1022	THR	2.8
1	A	53	PRO	2.8
1	A	674	ASP	2.8
1	A	51	LEU	2.8
3	C	21	GLN	2.8
1	A	700	PHE	2.8
3	C	149	MET	2.8
2	B	427	THR	2.6
1	A	94	PHE	2.6
3	C	33	TYR	2.6
1	A	592	VAL	2.6
2	B	662	LYS	2.6
1	A	544	ARG	2.6
1	A	91	ARG	2.5
2	B	424	THR	2.5
2	B	771	THR	2.5
3	C	30	LEU	2.4
1	A	508	THR	2.4
1	A	667	TRP	2.4
2	B	1048	PHE	2.3
1	A	122	LEU	2.3
1	A	510	ILE	2.3
1	A	593	PHE	2.3
2	B	348	LEU	2.3
2	B	346	GLU	2.2
1	A	431	PRO	2.2
2	B	449	ARG	2.2
1	A	536	THR	2.2
1	A	79	ARG	2.2
1	A	672	TYR	2.2
2	B	914	THR	2.2
3	C	98	GLN	2.1
1	A	507	GLN	2.1
1	A	675	MET	2.1
3	C	20	MET	2.0
1	A	760	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	103	GLU	2.0
2	B	478	HIS	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 monosaccharides in this entry.

6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	ZN	B	1101	1/1	0.91	0.16	94,94,94,94	0
5	ZN	A	801	1/1	0.96	0.12	95,95,95,95	0

6.5 Other polymers (i)

There are no such residues in this entry.