



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 2, 2017 – 11:11 AM EST

PDB ID : 1W6I
Title : plasmepsin II-pepstatin A complex
Authors : Lindberg, J.; Johansson, P.-O.; Rosenquist, A.; Kvarnstrom, I.; Vrang, L.; Samuelsson, B.; Unge, T.
Deposited on : 2004-08-18
Resolution : 2.70 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
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-20028442

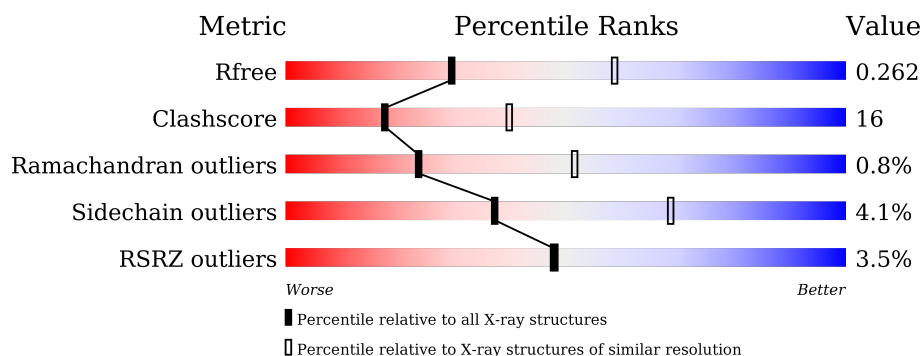
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	331	<div> <div>2%</div> <div> <div></div> <div>68%</div> <div>29%</div> <div>..</div> </div> </div>
1	C	331	<div> <div>5%</div> <div> <div></div> <div>67%</div> <div>29%</div> <div>..</div> </div> </div>
2	B	6	<div> <div></div> <div> <div>50%</div> <div>50%</div> </div> </div>
2	D	6	<div> <div></div> <div> <div>67%</div> <div>33%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5542 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 PLASMEPSIN 2 PRECURSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2606	1689	404	502	11			
1	C	329	Total	C	N	O	S	0	0	0
			2606	1689	404	502	11			

- Molecule 2 is a protein called PEPSTATIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	6	Total	C	N	O	0	0	0
			48	34	5	9			
2	D	6	Total	C	N	O	0	0	0
			48	34	5	9			

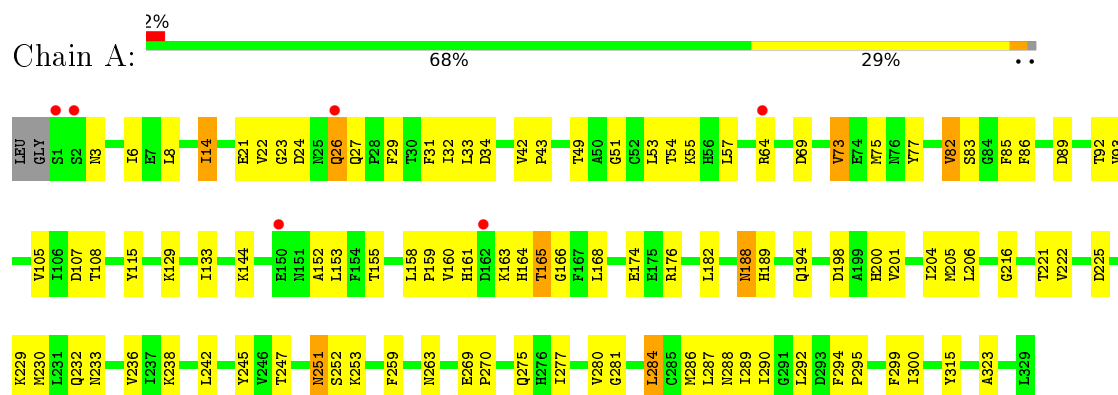
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	132	Total	O	0	0
			132	132		
3	C	100	Total	O	0	0
			100	100		
3	B	1	Total	O	0	0
			1	1		
3	D	1	Total	O	0	0
			1	1		

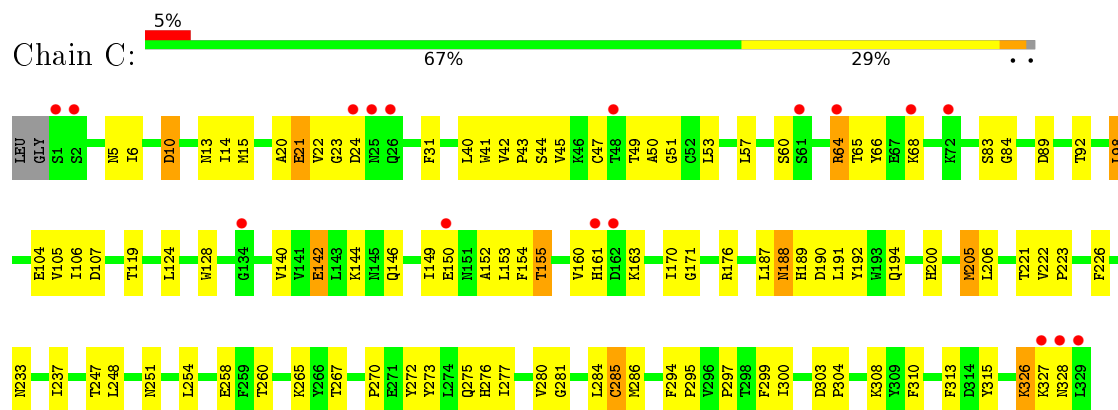
3 Residue-property plots

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: PLASMEPSIN 2 PRECURSOR



• Molecule 1: PLASMEPSIN 2 PRECURSOR



• Molecule 2: PEPSTATIN



• Molecule 2: PEPSTATIN



?	1330
V	1331
V	1332
?	1333
A	1334
?	1335

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	142.29 Å 142.29 Å 98.09 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 2.70 19.92 – 2.61	Depositor EDS
% Data completeness (in resolution range)	99.4 (25.00-2.70) 99.4 (19.92-2.61)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.64 (at 2.59 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.217 , 0.262 0.217 , 0.262	Depositor DCC
R_{free} test set	1590 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	33.0	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 57.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5542	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.83% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: STA, IVA

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	0.43	0/2673	0.69	1/3638 (0.0%)
1	C	0.41	0/2673	0.65	0/3638
2	B	0.77	0/17	2.52	1/21 (4.8%)
2	D	0.75	0/17	2.62	1/21 (4.8%)
All	All	0.42	0/5380	0.70	3/7318 (0.0%)

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
2	B	0	2
2	D	0	2
All	All	0	4

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	D	1331	VAL	CG1-CB-CG2	-8.14	97.87	110.90
2	B	1331	VAL	CG1-CB-CG2	-7.43	99.02	110.90
1	A	32	ILE	N-CA-C	-5.53	96.06	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	1333	STA	Mainchain,Peptide

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Mol	Chain	Res	Type	Group
2	D	1333	STA	Mainchain,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	2606	0	2536	83	0
1	C	2606	0	2536	84	0
2	B	48	0	60	1	0
2	D	48	0	60	0	0
3	A	132	0	0	3	0
3	B	1	0	0	1	0
3	C	100	0	0	1	0
3	D	1	0	0	0	0
All	All	5542	0	5192	167	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 16.

All (167) 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:221:THR:HB	1:A:300:ILE:HB	1.34	1.08
1:A:6:ILE:HD11	1:A:93:VAL:HG13	1.41	0.99
1:C:221:THR:HB	1:C:300:ILE:HB	1.41	0.99
1:A:236:VAL:HG22	1:A:247:THR:HG21	1.45	0.95
1:A:236:VAL:HG22	1:A:247:THR:CG2	2.03	0.89
1:C:44:SER:HB2	1:C:104:GLU:HG2	1.55	0.88
1:A:6:ILE:HD11	1:A:93:VAL:CG1	2.08	0.84
1:A:251:ASN:ND2	1:A:253:LYS:H	1.82	0.78
1:A:21:GLU:HG2	1:A:92:THR:HB	1.68	0.75
1:C:144:LYS:HA	1:C:149:ILE:HG22	1.69	0.75
1:C:49:THR:HG22	1:C:51:GLY:H	1.54	0.72
1:A:6:ILE:CD1	1:A:93:VAL:HG13	2.20	0.71
1:C:24:ASP:HB2	1:C:65:THR:OG1	1.91	0.71
1:A:290:ILE:HG13	2:B:1330:IVA:HG11	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:206:LEU:HD23	1:C:299:PHE:HE1	1.55	0.70
1:C:251:ASN:HB3	1:C:254:LEU:HG	1.73	0.69
1:C:160:VAL:HG11	1:C:163:LYS:HD3	1.74	0.69
1:A:188:ASN:HD21	1:A:194:GLN:HE21	1.42	0.68
1:C:23:GLY:HA2	1:C:89:ASP:OD1	1.94	0.68
1:C:276:HIS:HA	1:C:285:CYS:HB3	1.74	0.67
1:C:41:TRP:CE3	1:C:105:VAL:HG21	2.30	0.67
1:A:200:HIS:HB2	1:A:205:MET:HE2	1.78	0.66
1:A:287:LEU:O	1:A:289:ILE:HG13	1.97	0.65
1:A:251:ASN:C	1:A:251:ASN:HD22	1.98	0.65
1:C:260:THR:HG22	1:C:265:LYS:HG3	1.78	0.65
1:A:33:LEU:HD13	1:A:168:LEU:HD22	1.79	0.65
1:A:174:GLU:OE1	1:A:176:ARG:HD3	1.97	0.63
1:C:188:ASN:HB2	1:C:194:GLN:HG3	1.81	0.62
1:A:281:GLY:O	1:A:284:LEU:HB2	1.99	0.62
1:A:54:THR:HG22	1:A:54:THR:O	2.01	0.61
1:C:221:THR:HB	1:C:300:ILE:CB	2.24	0.61
1:C:277:ILE:HG22	1:C:280:VAL:HB	1.82	0.61
1:C:294:PHE:HB3	1:C:295:PRO:HD2	1.83	0.60
1:A:236:VAL:HA	1:A:247:THR:HG22	1.84	0.60
1:C:153:LEU:C	1:C:153:LEU:HD12	2.22	0.60
1:A:14:ILE:HD11	1:A:275:GLN:CD	2.23	0.59
1:C:223:PRO:HG2	1:C:226:PHE:HB2	1.84	0.59
1:C:14:ILE:CD1	1:C:275:GLN:HG3	2.33	0.59
1:C:163:LYS:HG2	1:C:327:LYS:HE3	1.85	0.59
1:C:45:VAL:HA	1:C:57:LEU:HD13	1.84	0.59
1:C:84:GLY:HA3	1:C:104:GLU:O	2.04	0.58
1:C:247:THR:HG22	1:C:248:LEU:N	2.17	0.58
1:C:222:VAL:HG23	1:C:299:PHE:CE2	2.39	0.57
1:C:273:TYR:HA	1:C:304:PRO:HB3	1.85	0.57
1:C:64:ARG:H	1:C:64:ARG:HD2	1.70	0.57
1:C:14:ILE:HD12	1:C:275:GLN:HG3	1.86	0.57
1:A:251:ASN:HD22	1:A:252:SER:N	2.03	0.56
1:C:281:GLY:HA3	1:C:284:LEU:HD22	1.87	0.56
1:C:47:CYS:HB2	1:C:106:ILE:HA	1.86	0.56
1:C:237:ILE:HD12	1:C:237:ILE:N	2.21	0.56
1:C:149:ILE:HG13	1:C:171:GLY:HA2	1.88	0.55
1:A:73:VAL:HG23	1:A:85:PHE:HA	1.89	0.55
1:A:69:ASP:HB3	1:A:86:PHE:O	2.08	0.54
1:A:188:ASN:C	1:A:188:ASN:HD22	2.11	0.54
1:A:277:ILE:HG13	1:A:280:VAL:CG2	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:LEU:C	1:A:153:LEU:HD12	2.28	0.54
1:A:236:VAL:HG13	1:A:247:THR:HG23	1.91	0.53
1:A:159:PRO:HB3	1:A:166:GLY:N	2.23	0.53
1:A:205:MET:HG2	1:A:206:LEU:N	2.22	0.53
1:C:200:HIS:CD2	1:C:205:MET:HB3	2.44	0.53
1:A:21:GLU:HB2	1:A:26:GLN:O	2.08	0.53
1:A:281:GLY:HA3	1:A:284:LEU:HD22	1.91	0.53
1:C:222:VAL:HG22	1:C:226:PHE:HB3	1.90	0.53
1:A:221:THR:HB	1:A:300:ILE:CB	2.23	0.52
1:A:163:LYS:HG2	1:A:164:HIS:CD2	2.45	0.52
1:C:277:ILE:HG13	1:C:286:MET:CE	2.40	0.52
1:C:326:LYS:HD3	1:C:328:ASN:HB2	1.91	0.52
1:A:198:ASP:O	1:A:259:PHE:HA	2.11	0.51
1:C:5:ASN:C	1:C:6:ILE:HD12	2.31	0.51
1:C:176:ARG:O	1:C:326:LYS:HD2	2.12	0.50
1:A:82:VAL:HG22	1:A:105:VAL:HG11	1.93	0.50
1:A:251:ASN:C	1:A:251:ASN:ND2	2.65	0.50
1:A:288:ASN:HA	3:A:2120:HOH:O	2.10	0.50
1:A:133:ILE:HD11	3:B:2133:HOH:O	2.10	0.50
1:A:294:PHE:HB3	1:A:295:PRO:CD	2.41	0.50
1:A:221:THR:HG23	1:A:292:LEU:HB3	1.92	0.50
1:C:152:ALA:O	1:C:315:TYR:HB2	2.12	0.50
1:A:269:GLU:HB2	1:A:270:PRO:HD2	1.94	0.50
1:A:53:LEU:HD23	1:A:53:LEU:C	2.33	0.50
1:C:297:PRO:HG2	3:C:2085:HOH:O	2.12	0.49
1:C:60:SER:HB2	1:C:66:TYR:CD1	2.48	0.49
1:C:23:GLY:HA2	1:C:89:ASP:CG	2.33	0.49
1:C:144:LYS:HE3	1:C:150:GLU:O	2.11	0.49
1:A:8:LEU:O	1:A:165:THR:HG22	2.12	0.49
1:A:34:ASP:OD1	1:A:216:GLY:HA3	2.13	0.48
1:A:269:GLU:H	1:A:269:GLU:CD	2.16	0.48
1:A:53:LEU:HD23	1:A:53:LEU:O	2.13	0.48
1:A:200:HIS:HB2	1:A:205:MET:CE	2.43	0.48
1:C:277:ILE:HG13	1:C:286:MET:HE3	1.95	0.48
1:C:128:TRP:CE3	1:C:190:ASP:HB3	2.48	0.48
1:A:49:THR:HG21	1:A:108:THR:OG1	2.13	0.47
1:A:83:SER:HB2	1:A:107:ASP:HB3	1.94	0.47
1:A:160:VAL:HB	1:A:163:LYS:HB3	1.96	0.47
1:C:222:VAL:HG23	1:C:299:PHE:CZ	2.49	0.47
1:A:75:MET:HB3	1:A:82:VAL:HG12	1.96	0.47
1:C:20:ALA:HB3	1:C:31:PHE:CE2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:LEU:O	1:C:98:LEU:HG	2.15	0.47
1:C:144:LYS:CA	1:C:149:ILE:HG22	2.41	0.47
1:A:251:ASN:HD22	1:A:253:LYS:H	1.56	0.47
1:C:270:PRO:HA	1:C:273:TYR:CZ	2.50	0.46
1:A:222:VAL:HG22	1:A:299:PHE:CE2	2.51	0.46
1:A:163:LYS:HG2	1:A:164:HIS:HD2	1.80	0.46
1:A:238:LYS:HE3	1:A:242:LEU:O	2.15	0.46
1:A:29:PHE:HB2	1:A:31:PHE:CE1	2.50	0.46
1:C:40:LEU:HD13	1:C:124:LEU:HG	1.98	0.46
1:C:142:GLU:HG3	1:C:146:GLN:HE21	1.81	0.46
1:A:152:ALA:O	1:A:315:TYR:HB2	2.16	0.46
1:C:15:MET:HE3	1:C:303:ASP:OD2	2.15	0.46
1:A:43:PRO:HG3	1:A:115:TYR:OH	2.16	0.45
1:C:6:ILE:N	1:C:6:ILE:HD12	2.30	0.45
1:C:160:VAL:CG1	1:C:163:LYS:HD3	2.46	0.45
1:C:50:ALA:HA	1:C:53:LEU:HD12	1.97	0.45
1:C:327:LYS:HG2	1:C:327:LYS:O	2.15	0.45
1:C:83:SER:HB2	1:C:107:ASP:HB3	1.98	0.45
1:A:188:ASN:HD21	1:A:194:GLN:NE2	2.11	0.45
1:A:158:LEU:HD23	1:A:158:LEU:N	2.32	0.44
1:A:75:MET:HG3	1:A:77:TYR:CE2	2.53	0.44
1:C:42:VAL:HG22	1:C:43:PRO:HD2	1.99	0.44
1:A:188:ASN:ND2	1:A:189:HIS:CD2	2.85	0.44
1:A:225:ASP:OD2	1:A:229:LYS:HE2	2.17	0.43
1:A:153:LEU:O	1:A:153:LEU:HD12	2.18	0.43
1:A:49:THR:HG22	1:A:51:GLY:H	1.83	0.43
1:C:140:VAL:HG21	1:C:154:PHE:CD2	2.53	0.43
1:A:42:VAL:HG13	1:A:43:PRO:HD2	1.99	0.43
1:C:222:VAL:HG21	1:C:226:PHE:CD1	2.53	0.43
1:A:245:TYR:O	1:A:286:MET:HA	2.19	0.43
1:A:294:PHE:HB3	1:A:295:PRO:HD2	1.99	0.43
1:C:42:VAL:CG2	1:C:43:PRO:HD2	2.48	0.43
1:C:270:PRO:HA	1:C:273:TYR:CE2	2.54	0.43
1:A:201:VAL:HG11	1:A:230:MET:HG3	2.01	0.43
1:A:188:ASN:ND2	1:A:194:GLN:HG3	2.33	0.43
1:C:163:LYS:HE2	1:C:327:LYS:HG3	2.00	0.43
1:C:258:GLU:HG2	1:C:267:THR:HG22	1.99	0.42
1:A:129:LYS:HB2	3:A:2054:HOH:O	2.19	0.42
1:A:24:ASP:OD1	1:A:64:ARG:N	2.53	0.42
1:C:10:ASP:OD1	1:C:13:ASN:HA	2.19	0.42
1:A:54:THR:O	1:A:54:THR:CG2	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:ILE:HD13	1:A:229:LYS:HB3	2.01	0.42
1:A:263:ASN:ND2	3:A:2106:HOH:O	2.50	0.42
1:C:149:ILE:HG12	1:C:170:ILE:HG22	2.01	0.42
1:C:222:VAL:CG2	1:C:223:PRO:HD2	2.50	0.42
1:C:22:VAL:HG12	1:C:23:GLY:N	2.35	0.42
1:C:21:GLU:HG2	1:C:92:THR:HB	2.02	0.42
1:C:189:HIS:HD2	1:C:194:GLN:NE2	2.17	0.42
1:C:144:LYS:HG2	1:C:152:ALA:HB2	2.02	0.41
1:A:182:LEU:HD23	1:A:323:ALA:HB2	2.02	0.41
1:A:290:ILE:HD13	1:A:290:ILE:HA	1.92	0.41
1:C:247:THR:HG22	1:C:248:LEU:O	2.21	0.41
1:C:153:LEU:HB2	1:C:313:PHE:O	2.21	0.41
1:C:155:THR:HG23	1:C:310:PHE:CE1	2.56	0.41
1:C:200:HIS:NE2	1:C:205:MET:HB3	2.35	0.41
1:C:275:GLN:O	1:C:285:CYS:HA	2.20	0.41
1:A:26:GLN:O	1:A:27:GLN:HB2	2.21	0.41
1:A:42:VAL:HA	1:A:43:PRO:HD3	1.96	0.41
1:C:272:TYR:CD1	1:C:308:LYS:HB2	2.55	0.41
1:C:327:LYS:O	1:C:327:LYS:CG	2.69	0.41
1:A:188:ASN:HD22	1:A:189:HIS:N	2.19	0.40
1:A:22:VAL:O	1:A:27:GLN:HB2	2.21	0.40
1:A:43:PRO:HB2	1:A:57:LEU:HD23	2.03	0.40
1:C:277:ILE:HG22	1:C:280:VAL:CB	2.49	0.40
1:C:277:ILE:HG22	1:C:280:VAL:CG2	2.51	0.40
1:A:55:LYS:HE3	1:A:115:TYR:CE1	2.56	0.40
1:C:124:LEU:HD23	1:C:124:LEU:C	2.42	0.40
1:A:23:GLY:HA2	1:A:89:ASP:CG	2.40	0.40
1:C:191:LEU:HB3	1:C:192:TYR:H	1.71	0.40
1:C:163:LYS:HB3	1:C:327:LYS:HE3	2.03	0.40
1:C:260:THR:CG2	1:C:265:LYS:HG3	2.48	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	327/331 (99%)	301 (92%)	22 (7%)	4 (1%)	16	39
1	C	327/331 (99%)	310 (95%)	16 (5%)	1 (0%)	46	75
2	B	3/6 (50%)	3 (100%)	0	0	100	100
2	D	3/6 (50%)	3 (100%)	0	0	100	100
All	All	660/674 (98%)	617 (94%)	38 (6%)	5 (1%)	24	51

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	161	HIS
1	C	161	HIS
1	A	232	GLN
1	A	14	ILE

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	294/295 (100%)	284 (97%)	10 (3%)	44	75
1	C	294/295 (100%)	280 (95%)	14 (5%)	31	62
2	B	2/2 (100%)	2 (100%)	0	100	100
2	D	2/2 (100%)	2 (100%)	0	100	100
All	All	592/594 (100%)	568 (96%)	24 (4%)	37	69

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	73	VAL
1	A	82	VAL

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Mol	Chain	Res	Type
1	A	144	LYS
1	A	155	THR
1	A	165	THR
1	A	188	ASN
1	A	233	ASN
1	A	251	ASN
1	A	284	LEU
1	C	10	ASP
1	C	21	GLU
1	C	64	ARG
1	C	68	LYS
1	C	98	LEU
1	C	119	THR
1	C	142	GLU
1	C	155	THR
1	C	187	LEU
1	C	188	ASN
1	C	205	MET
1	C	233	ASN
1	C	285	CYS
1	C	326	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	HIS
1	A	147	ASN
1	A	188	ASN
1	A	189	HIS
1	A	233	ASN
1	A	250	ASN
1	A	251	ASN
1	A	263	ASN
1	C	26	GLN
1	C	27	GLN
1	C	188	ASN
1	C	189	HIS
1	C	200	HIS
1	C	203	ASN
1	C	275	GLN
1	C	276	HIS
1	C	288	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	STA	B	1333	2	10,10,11	0.95	1 (10%)	10,12,14	1.60	3 (30%)
2	STA	B	1335	2	8,11,11	2.26	2 (25%)	8,14,14	1.51	2 (25%)
2	STA	D	1333	2	10,10,11	0.89	1 (10%)	10,12,14	1.67	3 (30%)
2	STA	D	1335	2	8,11,11	2.80	2 (25%)	8,14,14	1.51	2 (25%)

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	STA	B	1333	2	-	0/11/11/12	0/0/0/0
2	STA	B	1335	2	-	0/10/12/12	0/0/0/0
2	STA	D	1333	2	-	0/11/11/12	0/0/0/0
2	STA	D	1335	2	-	0/10/12/12	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1333	STA	CH-CA	2.34	1.55	1.53
2	B	1333	STA	CH-CA	2.59	1.56	1.53
2	B	1335	STA	OH-CH	3.75	1.51	1.43
2	D	1335	STA	OH-CH	4.12	1.52	1.43
2	B	1335	STA	CH-CA	4.94	1.58	1.53
2	D	1335	STA	CH-CA	6.31	1.60	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1333	STA	OH-CH-CM	-3.40	101.65	109.33
2	B	1333	STA	OH-CH-CM	-3.10	102.32	109.33
2	B	1333	STA	CB-CA-N	-2.92	100.81	108.96
2	D	1335	STA	CM-CH-CA	-2.73	108.52	112.68
2	D	1333	STA	CB-CA-N	-2.54	101.86	108.96
2	B	1335	STA	CM-CH-CA	-2.48	108.90	112.68
2	B	1333	STA	CG-CB-CA	2.01	120.43	115.75
2	D	1333	STA	CG-CB-CA	2.26	121.01	115.75
2	D	1335	STA	OH-CH-CM	2.71	114.36	109.61
2	B	1335	STA	OH-CH-CM	2.92	114.72	109.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	A	329/331 (99%)	-0.38	6 (1%) 71 72	11, 26, 45, 64	0
1	C	329/331 (99%)	-0.14	17 (5%) 31 30	14, 34, 59, 73	0
2	B	3/6 (50%)	-1.16	0 100 100	17, 17, 23, 23	0
2	D	3/6 (50%)	-0.88	0 100 100	24, 24, 27, 32	0
All	All	664/674 (98%)	-0.26	23 (3%) 48 48	11, 29, 54, 73	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	329	LEU	6.8
1	C	2	SER	5.4
1	A	1	SER	4.4
1	C	1	SER	3.6
1	C	64	ARG	3.1
1	C	48	THR	3.1
1	A	26	GLN	3.0
1	A	162	ASP	2.9
1	A	2	SER	2.9
1	C	134	GLY	2.7
1	A	64	ARG	2.7
1	C	72	LYS	2.7
1	C	161	HIS	2.6
1	C	162	ASP	2.5
1	C	150	GLU	2.5
1	C	68	LYS	2.4
1	C	24	ASP	2.4
1	C	26	GLN	2.4
1	C	61	SER	2.3
1	A	150	GLU	2.2
1	C	328	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	327	LYS	2.1
1	C	25	ASN	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	STA	D	1333	11/12	0.95	0.13	-	25,26,28,29	0
2	STA	B	1333	11/12	0.97	0.10	-	14,16,19,19	0
2	STA	B	1335	12/12	0.89	0.20	-	24,32,48,49	0
2	STA	D	1335	12/12	0.86	0.32	-	38,42,55,57	0

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.