Full wwPDB X-ray Structure Validation Report ⓘ

Dec 5, 2023 – 12:49 am GMT

PDB ID	:	1OCV
Title	:	the F116W mutant structure of ketosteroid isomerase from Comamonas testosteroni
Authors	:	Yun, Y.S.; Lee, T.-H.; Shin, S.
Deposited on	:	2003-02-11
Resolution	:	2.00 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 ⓘ 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 ⓘ) were used in the production of this report:

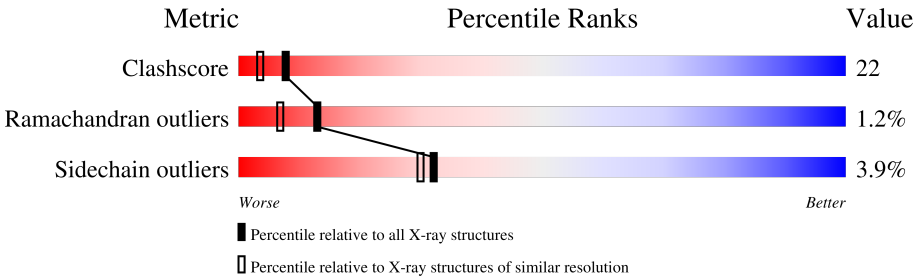
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.00 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))
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)

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%

Mol	Chain	Length	Quality of chain
1	A	125	75% 21%
1	B	125	72% 25%
1	C	125	66% 30%
1	D	125	67% 29%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4153 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 STEROID DELTA-ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	125	Total	C	N	O	S	0	0	0
			949	601	166	179	3			
1	B	125	Total	C	N	O	S	0	0	0
			949	601	166	179	3			
1	C	125	Total	C	N	O	S	0	0	0
			949	601	166	179	3			
1	D	125	Total	C	N	O	S	0	0	0
			949	601	166	179	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	83	ILE	THR	conflict	UNP P00947
B	283	ILE	THR	conflict	UNP P00947
C	483	ILE	THR	conflict	UNP P00947
D	683	ILE	THR	conflict	UNP P00947
A	116	TRP	PHE	engineered mutation	UNP P00947
B	316	TRP	PHE	engineered mutation	UNP P00947
C	516	TRP	PHE	engineered mutation	UNP P00947
D	716	TRP	PHE	engineered mutation	UNP P00947

- Molecule 2 is water.

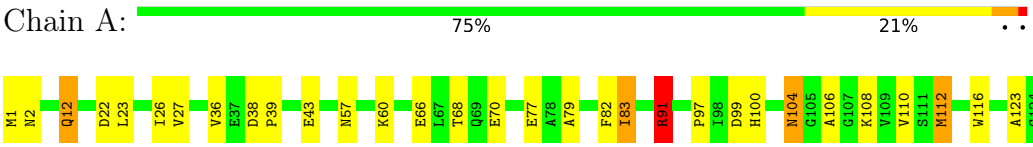
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	88	Total	O	0	0
			88	88		
2	B	86	Total	O	0	0
			86	86		
2	C	100	Total	O	0	0
			100	100		
2	D	83	Total	O	0	0
			83	83		

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. 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. 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.

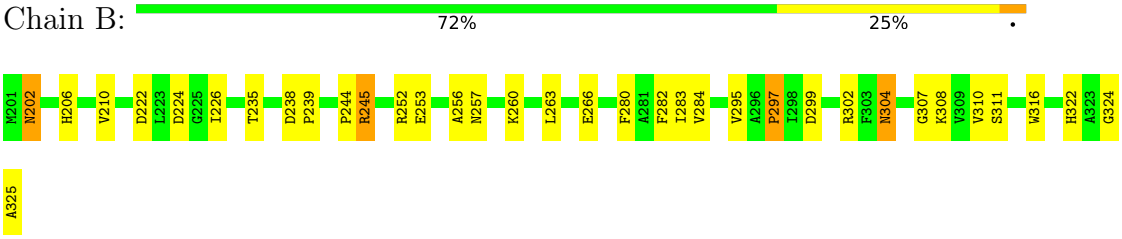
• Molecule 1: STEROID DELTA-ISOMERASE

Chain A:



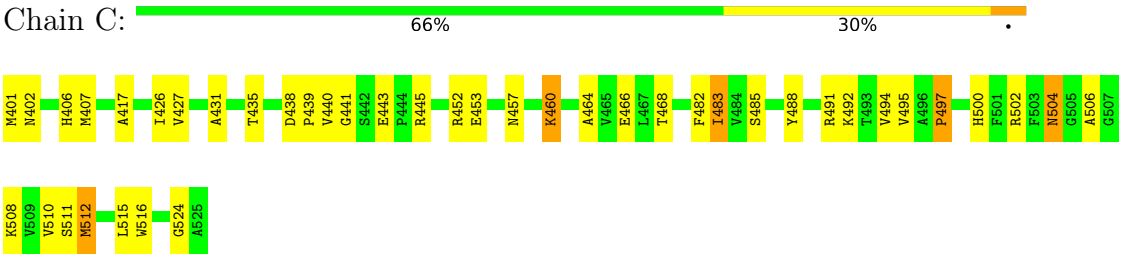
• Molecule 1: STEROID DELTA-ISOMERASE

Chain B:



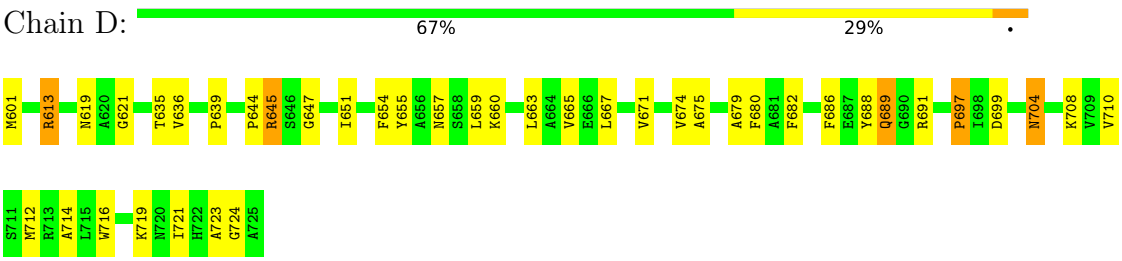
• Molecule 1: STEROID DELTA-ISOMERASE

Chain C:



• Molecule 1: STEROID DELTA-ISOMERASE

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α, β, γ	71.53Å 71.53Å 103.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.00 39.68 – 2.00	Depositor EDS
% Data completeness (in resolution range)	91.8 (50.00-2.00) 91.9 (39.68-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.55 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.229 , 0.282 0.243 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	19.1	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 38.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.39$, $< L^2 > = 0.21$	Xtriage
Estimated twinning fraction	0.104 for -h,-k,l 0.315 for h,-h-k,-l 0.105 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4153	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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

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	A	0.35	0/968	0.58	0/1314
1	B	0.36	0/968	0.58	0/1314
1	C	0.35	0/968	0.58	0/1314
1	D	0.34	0/968	0.58	0/1314
All	All	0.35	0/3872	0.58	0/5256

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

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	949	0	934	38	0
1	B	949	0	931	37	0
1	C	949	0	931	53	0
1	D	949	0	931	43	0
2	A	88	0	0	2	0
2	B	86	0	0	2	0
2	C	100	0	0	3	0
2	D	83	0	0	4	0
All	All	4153	0	3727	164	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 22.

All (164) 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:B:297:PRO:HB3	1:B:316:TRP:HB3	1.44	0.98
1:D:619:ASN:HD22	1:D:665:VAL:H	1.18	0.91
1:D:714:ALA:HB2	2:D:2036:HOH:O	1.75	0.87
1:A:66:GLU:HB2	1:A:83:ILE:HD11	1.58	0.85
1:C:497:PRO:HB3	1:C:516:TRP:HB3	1.58	0.85
1:A:97:PRO:HB3	1:A:116:TRP:HB3	1.60	0.81
1:B:304:ASN:HD21	1:B:308:LYS:H	1.30	0.79
1:B:304:ASN:HB3	1:B:310:VAL:HG11	1.66	0.77
1:B:304:ASN:ND2	1:B:308:LYS:H	1.83	0.77
1:C:483:ILE:H	1:C:483:ILE:HD13	1.51	0.75
1:D:636:VAL:HG13	1:D:712:MET:HE1	1.68	0.75
1:C:506:ALA:HB3	1:C:508:LYS:HE2	1.72	0.72
1:D:697:PRO:HB3	1:D:716:TRP:HB3	1.71	0.72
1:C:468:THR:HG23	1:C:483:ILE:HD12	1.72	0.71
1:B:282:PHE:CZ	1:B:297:PRO:HG2	2.26	0.70
1:B:283:ILE:HG13	2:B:2056:HOH:O	1.93	0.69
1:B:260:LYS:HD2	1:C:453:GLU:HG3	1.75	0.69
1:A:104:ASN:HB3	1:A:110:VAL:CG1	2.22	0.69
1:C:407:MET:HG2	2:C:2026:HOH:O	1.92	0.68
1:D:613:ARG:HH11	1:D:613:ARG:HG3	1.58	0.68
1:A:104:ASN:C	1:A:104:ASN:HD22	1.97	0.68
1:A:83:ILE:H	1:A:83:ILE:HD13	1.60	0.67
1:C:439:PRO:HB3	1:C:516:TRP:NE1	2.10	0.66
1:A:83:ILE:HG12	2:A:2060:HOH:O	1.95	0.65
1:D:619:ASN:ND2	1:D:665:VAL:H	1.94	0.65
1:C:504:ASN:HB3	1:C:510:VAL:HG11	1.76	0.65
1:C:504:ASN:C	1:C:504:ASN:HD22	2.00	0.65
1:D:704:ASN:C	1:D:704:ASN:HD22	1.99	0.65
1:D:704:ASN:HB3	1:D:710:VAL:HG11	1.79	0.64
1:A:104:ASN:HB3	1:A:110:VAL:HG11	1.80	0.64
1:C:466:GLU:HB2	1:C:483:ILE:HD11	1.79	0.64
1:D:682:PHE:CZ	1:D:697:PRO:HG2	2.33	0.63
1:A:43:GLU:H	1:A:43:GLU:CD	2.02	0.63
1:C:483:ILE:HD13	1:C:483:ILE:N	2.13	0.63
1:C:502:ARG:HB3	1:C:511:SER:HB3	1.79	0.62
1:A:39:PRO:HB3	1:A:116:TRP:NE1	2.14	0.62
1:D:655:TYR:HB2	2:D:2045:HOH:O	1.99	0.62
1:C:482:PHE:CZ	1:C:497:PRO:HG2	2.35	0.62
1:A:66:GLU:O	1:A:83:ILE:HD13	2.01	0.61
1:C:468:THR:CG2	1:C:483:ILE:HD12	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:PRO:HB3	1:A:116:TRP:CE2	2.36	0.60
1:A:104:ASN:HD21	1:A:108:LYS:H	1.49	0.60
1:D:636:VAL:HG13	1:D:712:MET:CE	2.31	0.60
1:C:504:ASN:HB3	1:C:510:VAL:CG1	2.31	0.60
1:A:1:MET:HE3	1:A:2:ASN:H	1.67	0.59
1:B:224:ASP:OD2	1:B:252:ARG:NH2	2.34	0.59
1:D:645:ARG:HH11	1:D:645:ARG:HG2	1.68	0.59
1:B:304:ASN:HB3	1:B:310:VAL:CG1	2.32	0.59
1:B:222:ASP:O	1:B:226:ILE:HG13	2.03	0.58
1:A:104:ASN:ND2	1:A:108:LYS:H	2.02	0.58
1:A:23:LEU:O	1:A:27:VAL:HG23	2.04	0.58
1:D:704:ASN:HB3	1:D:710:VAL:CG1	2.33	0.58
1:D:719:LYS:HB3	1:D:719:LYS:NZ	2.18	0.58
1:A:82:PHE:CZ	1:A:97:PRO:HG2	2.38	0.58
1:C:485:SER:HB3	1:C:492:LYS:NZ	2.18	0.58
1:D:688:TYR:O	1:D:689:GLN:HB2	2.04	0.57
1:A:83:ILE:HD13	1:A:83:ILE:N	2.20	0.57
1:C:466:GLU:O	1:C:483:ILE:HD13	2.04	0.57
1:B:297:PRO:CB	1:B:316:TRP:HB3	2.28	0.56
1:D:724:GLY:HA3	2:D:2082:HOH:O	2.05	0.56
1:C:497:PRO:CB	1:C:516:TRP:HB3	2.32	0.56
1:B:302:ARG:HB2	1:B:311:SER:HB2	1.88	0.56
1:B:302:ARG:HD3	1:B:311:SER:HB2	1.86	0.56
1:B:295:VAL:HG12	1:B:297:PRO:HD3	1.86	0.56
1:D:645:ARG:N	1:D:645:ARG:HD2	2.20	0.55
1:D:704:ASN:ND2	1:D:708:LYS:H	2.03	0.55
1:C:485:SER:HB3	1:C:492:LYS:HZ1	1.72	0.55
1:C:483:ILE:HG12	2:C:2069:HOH:O	2.08	0.54
1:C:406:HIS:HD2	2:C:2026:HOH:O	1.91	0.53
1:D:704:ASN:HD21	1:D:708:LYS:H	1.55	0.53
1:A:57:ASN:HB3	2:A:2053:HOH:O	2.08	0.53
1:B:266:GLU:OE1	1:B:283:ILE:HD11	2.08	0.53
1:B:302:ARG:O	1:B:310:VAL:HG22	2.09	0.53
1:B:257:ASN:HA	1:B:260:LYS:HZ3	1.73	0.53
1:D:645:ARG:NH1	1:D:654:PHE:CD1	2.77	0.53
1:A:104:ASN:OD1	1:A:108:LYS:HE3	2.10	0.52
1:D:613:ARG:HG3	1:D:613:ARG:NH1	2.24	0.52
1:D:635:THR:CG2	1:D:644:PRO:HB2	2.39	0.52
1:B:324:GLY:O	1:B:325:ALA:HB3	2.10	0.52
1:C:439:PRO:HB3	1:C:516:TRP:CE2	2.46	0.51
1:C:435:THR:HA	1:C:445:ARG:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:492:LYS:HE3	1:C:494:VAL:CG2	2.41	0.51
1:B:304:ASN:C	1:B:304:ASN:HD22	2.14	0.50
1:B:257:ASN:O	1:B:260:LYS:HG2	2.11	0.50
1:C:427:VAL:HG21	1:C:452:ARG:HG3	1.94	0.50
1:A:106:ALA:HB3	1:A:108:LYS:HE2	1.93	0.50
1:C:417:ALA:HB1	1:C:426:ILE:HG13	1.94	0.49
1:D:645:ARG:NH1	1:D:654:PHE:HD1	2.10	0.49
1:A:22:ASP:O	1:A:26:ILE:HG13	2.13	0.49
1:C:431:ALA:HA	1:C:508:LYS:HD2	1.95	0.49
1:A:91:ARG:CZ	1:A:123:ALA:HB1	2.43	0.48
1:D:704:ASN:C	1:D:704:ASN:ND2	2.65	0.48
1:A:97:PRO:CB	1:A:116:TRP:HB3	2.38	0.48
1:C:502:ARG:HG3	1:C:502:ARG:HH11	1.78	0.48
1:A:104:ASN:C	1:A:104:ASN:ND2	2.66	0.48
1:C:440:VAL:HG12	1:D:675:ALA:HB3	1.94	0.48
1:C:504:ASN:ND2	1:C:508:LYS:H	2.12	0.48
1:D:688:TYR:O	1:D:689:GLN:CB	2.62	0.48
1:C:401:MET:HE3	1:C:402:ASN:H	1.78	0.48
1:B:297:PRO:HB3	1:B:316:TRP:CB	2.30	0.47
1:D:657:ASN:HA	1:D:660:LYS:NZ	2.29	0.47
1:D:639:PRO:HB3	1:D:716:TRP:CE2	2.50	0.47
1:A:77:GLU:OE1	1:A:100:HIS:CE1	2.68	0.47
1:C:497:PRO:HB3	1:C:516:TRP:CB	2.38	0.46
1:C:500:HIS:O	1:C:512:MET:HA	2.16	0.46
1:B:257:ASN:HA	1:B:260:LYS:NZ	2.30	0.46
1:D:621:GLY:HA2	1:D:659:LEU:HD22	1.97	0.46
1:B:263:LEU:HD22	1:B:284:VAL:CG1	2.45	0.46
1:B:302:ARG:HD3	1:B:311:SER:CB	2.46	0.46
1:C:492:LYS:HD3	1:C:492:LYS:C	2.36	0.46
1:A:57:ASN:O	1:A:60:LYS:HD3	2.15	0.45
1:B:245:ARG:N	1:B:245:ARG:HD2	2.31	0.45
1:C:504:ASN:OD1	1:C:508:LYS:HE3	2.16	0.45
1:A:38:ASP:HA	1:A:39:PRO:HA	1.77	0.45
1:D:686:PHE:HE1	1:D:688:TYR:HB2	1.81	0.45
1:C:443:GLU:OE2	1:C:443:GLU:N	2.36	0.45
1:D:601:MET:HB2	2:D:2014:HOH:O	2.17	0.45
1:D:657:ASN:O	1:D:660:LYS:HG2	2.16	0.45
1:B:235:THR:HG23	1:B:244:PRO:HB2	1.99	0.45
1:B:280:PHE:CZ	1:B:299:ASP:HB2	2.52	0.45
1:C:512:MET:SD	1:C:512:MET:C	2.95	0.45
1:D:686:PHE:CE1	1:D:688:TYR:HB2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:671:VAL:HG22	1:D:680:PHE:HB3	1.99	0.44
1:C:464:ALA:HB3	1:C:485:SER:HB2	1.99	0.44
1:C:504:ASN:HD21	1:C:508:LYS:H	1.65	0.44
1:A:104:ASN:HB3	1:A:110:VAL:HG13	1.96	0.44
1:C:483:ILE:N	1:C:483:ILE:CD1	2.80	0.43
1:C:508:LYS:HE3	1:C:508:LYS:HB2	1.80	0.43
1:C:495:VAL:HG12	1:C:497:PRO:HD3	1.99	0.43
1:C:488:TYR:O	1:C:491:ARG:HB3	2.18	0.43
1:D:647:GLY:O	1:D:651:ILE:HG13	2.19	0.43
1:A:68:THR:CG2	1:A:83:ILE:HD12	2.49	0.43
1:C:504:ASN:C	1:C:504:ASN:ND2	2.69	0.43
1:C:492:LYS:HD2	1:C:524:GLY:O	2.19	0.43
1:A:12:GLN:HE21	1:A:12:GLN:HB2	1.55	0.42
1:D:679:ALA:HA	1:D:699:ASP:O	2.19	0.42
1:A:70:GLU:OE2	1:B:322:HIS:HE1	2.02	0.42
1:A:79:ALA:HA	1:A:99:ASP:O	2.19	0.42
1:A:97:PRO:HB3	1:A:116:TRP:CB	2.42	0.42
1:C:502:ARG:HG3	1:C:502:ARG:NH1	2.34	0.42
1:B:238:ASP:HA	1:B:239:PRO:HA	1.85	0.42
1:C:457:ASN:HA	1:C:460:LYS:CE	2.49	0.42
1:A:91:ARG:HG3	1:A:91:ARG:HH11	1.85	0.42
1:B:235:THR:CG2	1:B:244:PRO:HB2	2.50	0.41
1:B:257:ASN:HD22	1:B:260:LYS:NZ	2.18	0.41
1:D:657:ASN:HA	1:D:660:LYS:HZ3	1.85	0.41
1:D:682:PHE:CE2	1:D:697:PRO:HG2	2.55	0.41
1:C:438:ASP:HA	1:C:439:PRO:HA	1.80	0.41
1:C:441:GLY:HA3	1:D:675:ALA:HB2	2.02	0.41
1:A:66:GLU:O	1:A:82:PHE:HB2	2.21	0.41
1:D:663:LEU:CD2	1:D:686:PHE:HB3	2.51	0.41
1:A:108:LYS:HE3	1:A:108:LYS:HB2	1.85	0.41
1:B:253:GLU:HG3	1:C:460:LYS:HG3	2.02	0.41
1:A:1:MET:HG3	1:A:2:ASN:N	2.36	0.41
1:C:515:LEU:CD2	1:D:674:VAL:HG12	2.50	0.41
1:A:36:VAL:HA	1:A:112:MET:HG3	2.02	0.41
1:B:256:ALA:HB3	1:C:460:LYS:HD3	2.02	0.41
1:C:492:LYS:HE3	1:C:494:VAL:HG23	2.02	0.41
1:D:691:ARG:HH21	1:D:723:ALA:CB	2.34	0.41
1:D:716:TRP:HZ3	1:D:721:ILE:HD11	1.85	0.41
1:B:206:HIS:O	1:B:210:VAL:HG23	2.21	0.40
1:B:238:ASP:OD1	1:B:238:ASP:C	2.60	0.40
1:B:304:ASN:ND2	1:B:307:GLY:H	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:ASN:HA	2:B:2017:HOH:O	2.21	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	123/125 (98%)	115 (94%)	7 (6%)	1 (1%)	19 13
1	B	123/125 (98%)	119 (97%)	2 (2%)	2 (2%)	9 4
1	C	123/125 (98%)	120 (98%)	2 (2%)	1 (1%)	19 13
1	D	123/125 (98%)	117 (95%)	4 (3%)	2 (2%)	9 4
All	All	492/500 (98%)	471 (96%)	15 (3%)	6 (1%)	13 7

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	689	GLN
1	A	91	ARG
1	B	202	ASN
1	B	297	PRO
1	D	697	PRO
1	C	497	PRO

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	95/95 (100%)	90 (95%)	5 (5%)	22	18
1	B	95/95 (100%)	93 (98%)	2 (2%)	53	57
1	C	95/95 (100%)	91 (96%)	4 (4%)	30	27
1	D	95/95 (100%)	91 (96%)	4 (4%)	30	27
All	All	380/380 (100%)	365 (96%)	15 (4%)	32	30

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	83	ILE
1	A	91	ARG
1	A	104	ASN
1	A	112	MET
1	B	245	ARG
1	B	304	ASN
1	C	460	LYS
1	C	483	ILE
1	C	504	ASN
1	C	512	MET
1	D	613	ARG
1	D	645	ARG
1	D	667	LEU
1	D	704	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	57	ASN
1	A	89	GLN
1	A	100	HIS
1	A	104	ASN
1	B	212	GLN
1	B	257	ASN
1	B	289	GLN
1	B	304	ASN
1	B	322	HIS
1	C	412	GLN

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Mol	Chain	Res	Type
1	C	489	GLN
1	C	504	ASN
1	C	522	HIS
1	D	612	GLN
1	D	619	ASN
1	D	657	ASN
1	D	704	ASN
1	D	722	HIS

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

There are no monosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.