Full wwPDB X-ray Structure Validation Report (i)

Jun 18, 2024 – 11:22 PM EDT

PDB ID	:	4G04
Title	:	High-resolution Crystal Structural Variance Analysis between Recombinant and Wild-type Human Serum Albumin
Authors	:	Cao, H.L.; Yin, D.C.
Deposited on	:	2012-07-09
Resolution	:	2.30 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:

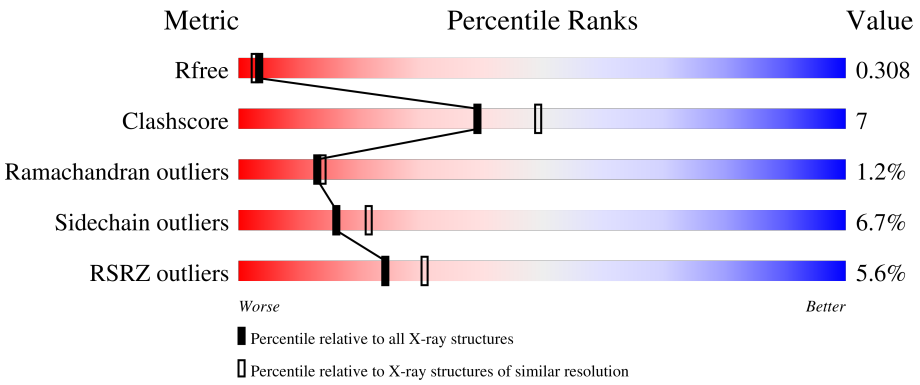
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	585	5% 76% 21% • •
1	B	585	6% 76% 20% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 9259 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 Serum albumin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	578	4599	2903	776	879	41	0	0	0
1	B	578	4599	2903	776	879	41	0	0	0

- Molecule 2 is water.

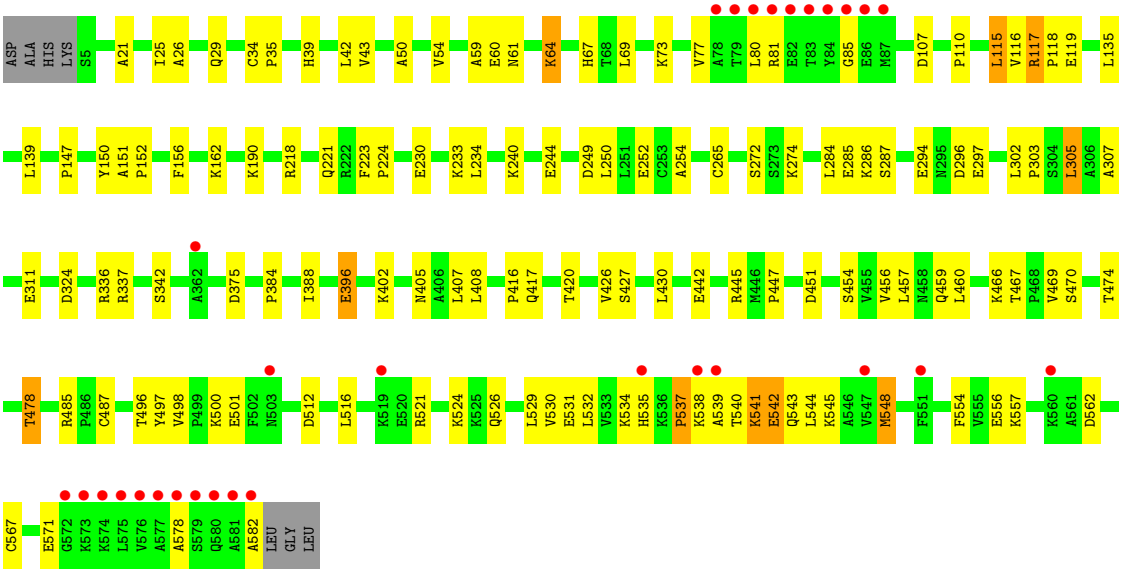
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	29	29	29	0	0
2	B	32	32	32	0	0

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.

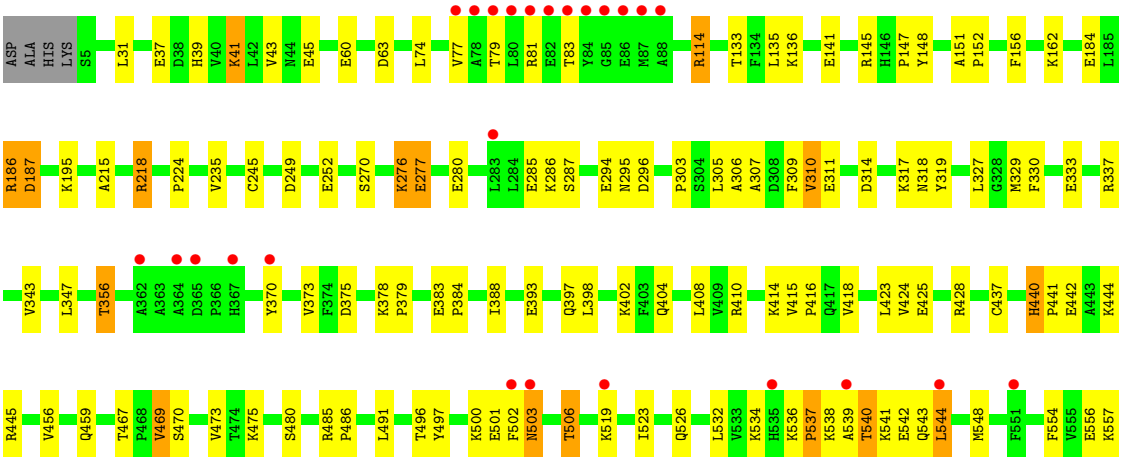
• Molecule 1: Serum albumin

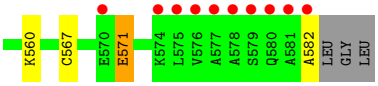
Chain A:



• Molecule 1: Serum albumin

Chain B:




K560
C567
E570
E571
K574
L575
V576
A577
A578
S579
Q580
A581
A582
LEU
GLY
LEU

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	58.74Å 58.66Å 95.73Å 87.85° 75.92° 73.47°	Depositor
Resolution (Å)	47.37 – 2.30 47.37 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.8 (47.37-2.30) 82.7 (47.37-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$< I/\sigma(I) >$ [1]	1.49 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.240 , 0.317 0.234 , 0.308	Depositor DCC
R_{free} test set	1998 reflections (4.20%)	wwPDB-VP
Wilson B-factor (Å^2)	39.2	Xtriage
Anisotropy	0.248	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 41.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	9259	wwPDB-VP
Average B, all atoms (Å^2)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.02% 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)

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	0.44	0/4688	0.57	0/6324
1	B	0.47	0/4688	0.57	0/6324
All	All	0.45	0/9376	0.57	0/12648

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	4599	0	4518	69	0
1	B	4599	0	4518	65	0
2	A	29	0	0	0	0
2	B	32	0	0	1	0
All	All	9259	0	9036	134	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 7.

All (134) 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:303:PRO:O	1:A:337:ARG:NH1	2.11	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:303:PRO:O	1:B:337:ARG:NH1	2.11	0.82
1:B:224:PRO:HD2	1:B:296:ASP:HB3	1.68	0.75
1:A:474:THR:O	1:A:478:THR:OG1	2.07	0.72
1:A:516:LEU:O	1:A:521:ARG:NH2	2.23	0.70
1:B:540:THR:HA	1:B:543:GLN:HE21	1.58	0.69
1:A:537:PRO:O	1:A:539:ALA:N	2.24	0.69
1:A:531:GLU:O	1:A:535:HIS:ND1	2.26	0.68
1:B:114:ARG:HH11	1:B:114:ARG:HB3	1.59	0.67
1:A:578:ALA:HA	1:A:582:ALA:HB3	1.77	0.66
1:A:224:PRO:HD2	1:A:296:ASP:HB3	1.79	0.65
1:B:384:PRO:O	1:B:388:ILE:HG12	1.99	0.63
1:B:437:CYS:O	1:B:440:HIS:HB2	2.00	0.61
1:B:156:PHE:HE1	1:B:285:GLU:HG3	1.64	0.61
1:B:414:LYS:HD3	1:B:491:LEU:HB2	1.84	0.60
1:A:427:SER:HA	1:A:430:LEU:HD12	1.84	0.59
1:B:156:PHE:CE1	1:B:285:GLU:HG3	2.37	0.59
1:A:305:LEU:HD11	1:A:337:ARG:HD2	1.85	0.59
1:A:540:THR:HA	1:A:543:GLN:HE21	1.67	0.59
1:A:156:PHE:HE1	1:A:285:GLU:HG3	1.68	0.58
1:B:416:PRO:O	1:B:534:LYS:HE2	2.04	0.58
1:B:306:ALA:HA	1:B:310:VAL:HG23	1.86	0.57
1:B:329:MET:O	1:B:333:GLU:HG2	2.04	0.57
1:A:384:PRO:O	1:A:388:ILE:HG12	2.05	0.56
1:A:296:ASP:OD1	1:A:297:GLU:N	2.33	0.56
1:A:42:LEU:HD22	1:A:73:LYS:HG3	1.88	0.56
1:A:156:PHE:CE1	1:A:285:GLU:HG3	2.41	0.56
1:B:314:ASP:HB3	1:B:317:LYS:HB3	1.87	0.56
1:B:276:LYS:O	1:B:280:GLU:HG2	2.05	0.56
1:B:195:LYS:NZ	2:B:631:HOH:O	2.40	0.55
1:B:519:LYS:O	1:B:523:ILE:HG13	2.06	0.55
1:B:141:GLU:OE1	1:B:145:ARG:NH1	2.30	0.54
1:B:276:LYS:HG3	1:B:277:GLU:N	2.22	0.54
1:A:61:ASN:HB3	1:A:64:LYS:HE2	1.90	0.53
1:A:408:LEU:HD11	1:A:526:GLN:HB3	1.90	0.52
1:A:554:PHE:CE1	1:A:571:GLU:HB2	2.44	0.52
1:A:223:PHE:CD1	1:A:272:SER:HB2	2.46	0.51
1:B:393:GLU:HG2	1:B:397:GLN:NE2	2.25	0.51
1:B:567:CYS:O	1:B:571:GLU:HG2	2.10	0.51
1:B:415:VAL:HG11	1:B:473:VAL:HG23	1.91	0.51
1:B:398:LEU:HB3	1:B:402:LYS:HB2	1.93	0.51
1:B:428:ARG:NE	1:B:526:GLN:OE1	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:PRO:O	1:A:534:LYS:HE2	2.12	0.50
1:B:356:THR:HG21	1:B:373:VAL:HG22	1.94	0.50
1:B:503:ASN:HB2	1:B:506:THR:OG1	2.12	0.50
1:A:240:LYS:HE2	1:A:244:GLU:OE2	2.11	0.50
1:B:186:ARG:NH1	1:B:187:ASP:OD1	2.45	0.50
1:B:218:ARG:NH2	1:B:295:ASN:OD1	2.45	0.50
1:A:135:LEU:HD11	1:A:162:LYS:HG3	1.94	0.49
1:A:223:PHE:HD1	1:A:272:SER:HB2	1.75	0.49
1:A:408:LEU:HD12	1:A:529:LEU:HD23	1.94	0.49
1:A:29:GLN:HG2	1:A:147:PRO:HA	1.94	0.49
1:A:61:ASN:O	1:A:64:LYS:HB2	2.13	0.49
1:A:152:PRO:HG3	1:A:254:ALA:HA	1.95	0.48
1:B:410:ARG:O	1:B:414:LYS:HG3	2.14	0.48
1:A:21:ALA:O	1:A:25:ILE:HG13	2.14	0.48
1:A:26:ALA:HB2	1:A:250:LEU:HD12	1.95	0.48
1:A:64:LYS:HB3	1:A:69:LEU:HG	1.94	0.48
1:A:405:ASN:ND2	1:A:548:MET:SD	2.87	0.47
1:B:404:GLN:HG2	1:B:428:ARG:HA	1.96	0.47
1:A:537:PRO:C	1:A:539:ALA:H	2.18	0.47
1:A:50:ALA:O	1:A:54:VAL:HG23	2.14	0.47
1:B:135:LEU:HD11	1:B:162:LYS:HG3	1.96	0.47
1:B:441:PRO:HG2	1:B:444:LYS:HD3	1.96	0.46
1:A:249:ASP:HB3	1:A:252:GLU:CG	2.44	0.46
1:A:426:VAL:HG21	1:A:460:LEU:HB2	1.96	0.46
1:A:265:CYS:SG	1:A:286:LYS:HD2	2.56	0.46
1:B:456:VAL:O	1:B:459:GLN:HB3	2.16	0.46
1:A:540:THR:HA	1:A:543:GLN:NE2	2.30	0.45
1:A:420:THR:OG1	1:A:531:GLU:OE2	2.21	0.45
1:B:538:LYS:HB2	1:B:539:ALA:H	1.45	0.45
1:A:302:LEU:HB3	1:A:337:ARG:NH1	2.32	0.45
1:A:416:PRO:HB2	1:A:497:TYR:CE1	2.52	0.45
1:B:383:GLU:HB3	1:B:384:PRO:HD3	1.99	0.45
1:A:218:ARG:HH21	1:A:221:GLN:HB2	1.82	0.44
1:B:63:ASP:OD1	1:B:63:ASP:N	2.39	0.44
1:B:485:ARG:HB3	1:B:486:PRO:HD3	1.99	0.44
1:A:67:HIS:NE2	1:A:249:ASP:OD1	2.51	0.44
1:A:117:ARG:HA	1:A:118:PRO:HD3	1.81	0.44
1:B:41:LYS:O	1:B:45:GLU:HG3	2.18	0.44
1:B:39:HIS:O	1:B:43:VAL:HG23	2.18	0.43
1:B:310:VAL:HG12	1:B:370:TYR:O	2.18	0.43
1:A:342:SER:HA	1:A:447:PRO:HA	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:408:LEU:HD11	1:B:526:GLN:HB3	1.99	0.43
1:A:416:PRO:HB2	1:A:497:TYR:CD1	2.53	0.43
1:B:343:VAL:O	1:B:347:LEU:HG	2.17	0.43
1:A:230:GLU:O	1:A:234:LEU:HG	2.19	0.43
1:A:402:LYS:HD3	1:A:545:LYS:HE3	2.00	0.43
1:B:133:THR:HA	1:B:136:LYS:HE3	2.00	0.43
1:B:307:ALA:HA	1:B:311:GLU:HB2	1.99	0.43
1:A:151:ALA:HB3	1:A:152:PRO:HD3	2.01	0.43
1:B:309:PHE:CZ	1:B:330:PHE:HA	2.54	0.43
1:A:540:THR:HG21	1:A:544:LEU:HD21	2.01	0.43
1:B:428:ARG:NE	1:B:523:ILE:HG12	2.34	0.43
1:B:424:VAL:HG12	1:B:428:ARG:HD2	2.01	0.43
1:A:39:HIS:O	1:A:43:VAL:HG23	2.18	0.42
1:B:151:ALA:HB3	1:B:152:PRO:HD3	1.99	0.42
1:B:536:LYS:HE3	1:B:582:ALA:O	2.20	0.42
1:A:107:ASP:O	1:A:110:PRO:HD3	2.19	0.42
1:B:249:ASP:HB3	1:B:252:GLU:CG	2.50	0.42
1:B:501:GLU:HG3	1:B:502:PHE:H	1.84	0.42
1:A:396:GLU:H	1:A:396:GLU:HG2	1.70	0.42
1:A:567:CYS:O	1:A:571:GLU:HG2	2.20	0.42
1:A:139:LEU:HA	1:A:139:LEU:HD23	1.76	0.42
1:A:417:GLN:HB3	1:A:469:VAL:HG12	2.02	0.42
1:B:319:TYR:HE1	1:B:327:LEU:HD11	1.84	0.42
1:B:393:GLU:HG2	1:B:397:GLN:HE21	1.85	0.42
1:B:467:THR:O	1:B:469:VAL:N	2.53	0.42
1:B:537:PRO:HB2	1:B:538:LYS:H	1.62	0.42
1:B:541:LYS:HB2	1:B:542:GLU:H	1.73	0.42
1:A:274:LYS:HE2	1:A:296:ASP:HA	2.01	0.41
1:A:457:LEU:HD23	1:A:457:LEU:HA	1.84	0.41
1:A:485:ARG:O	1:A:485:ARG:HD2	2.20	0.41
1:A:541:LYS:HB3	1:A:542:GLU:H	1.54	0.41
1:B:31:LEU:HG	1:B:74:LEU:HD22	2.02	0.41
1:B:215:ALA:HB3	1:B:235:VAL:HG13	2.02	0.41
1:A:307:ALA:HA	1:A:311:GLU:HB2	2.02	0.41
1:A:451:ASP:O	1:A:454:SER:HB2	2.20	0.41
1:B:416:PRO:HB2	1:B:497:TYR:CD1	2.55	0.41
1:B:544:LEU:HD22	1:B:544:LEU:HA	1.77	0.41
1:B:548:MET:HE3	1:B:548:MET:HA	2.02	0.41
1:B:428:ARG:CZ	1:B:523:ILE:HG12	2.51	0.41
1:A:81:ARG:HG3	1:A:85:GLY:O	2.21	0.41
1:A:284:LEU:HD23	1:A:284:LEU:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:407:LEU:HD13	1:A:430:LEU:HB3	2.02	0.41
1:A:456:VAL:O	1:A:459:GLN:HB3	2.21	0.41
1:A:526:GLN:O	1:A:530:VAL:HG23	2.21	0.41
1:B:418:VAL:HG12	1:B:423:LEU:HG	2.02	0.40
1:B:378:LYS:HB3	1:B:379:PRO:HD3	2.03	0.40
1:A:34:CYS:HA	1:A:35:PRO:HD3	1.86	0.40
1:A:115:LEU:HD13	1:A:115:LEU:HA	1.87	0.40
1:A:442:GLU:HA	1:A:445:ARG:HB2	2.02	0.40
1:B:442:GLU:HG2	1:B:445:ARG:HD2	2.03	0.40
1:B:147:PRO:HG2	1:B:148:TYR:CD1	2.57	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	576/585 (98%)	543 (94%)	24 (4%)	9 (2%)	9 9
1	B	576/585 (98%)	538 (93%)	33 (6%)	5 (1%)	17 20
All	All	1152/1170 (98%)	1081 (94%)	57 (5%)	14 (1%)	13 14

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	GLU
1	A	537	PRO
1	A	538	LYS
1	B	276	LYS
1	A	501	GLU
1	A	541	LYS
1	B	60	GLU
1	B	537	PRO

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Mol	Chain	Res	Type
1	A	80	LEU
1	B	83	THR
1	B	560	LYS
1	A	59	ALA
1	A	466	LYS
1	A	150	TYR

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	506/511 (99%)	476 (94%)	30 (6%)	19 27
1	B	506/511 (99%)	468 (92%)	38 (8%)	13 17
All	All	1012/1022 (99%)	944 (93%)	68 (7%)	16 21

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	LYS
1	A	77	VAL
1	A	115	LEU
1	A	116	VAL
1	A	117	ARG
1	A	119	GLU
1	A	190	LYS
1	A	233	LYS
1	A	287	SER
1	A	294	GLU
1	A	305	LEU
1	A	324	ASP
1	A	336	ARG
1	A	375	ASP
1	A	396	GLU
1	A	467	THR
1	A	470	SER

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Mol	Chain	Res	Type
1	A	478	THR
1	A	487	CYS
1	A	496	THR
1	A	498	VAL
1	A	500	LYS
1	A	512	ASP
1	A	524	LYS
1	A	532	LEU
1	A	542	GLU
1	A	548	MET
1	A	556	GLU
1	A	557	LYS
1	A	562	ASP
1	B	37	GLU
1	B	41	LYS
1	B	77	VAL
1	B	79	THR
1	B	81	ARG
1	B	114	ARG
1	B	184	GLU
1	B	186	ARG
1	B	187	ASP
1	B	218	ARG
1	B	245	CYS
1	B	270	SER
1	B	277	GLU
1	B	286	LYS
1	B	287	SER
1	B	294	GLU
1	B	305	LEU
1	B	310	VAL
1	B	318	ASN
1	B	356	THR
1	B	375	ASP
1	B	425	GLU
1	B	440	HIS
1	B	469	VAL
1	B	470	SER
1	B	475	LYS
1	B	480	SER
1	B	496	THR
1	B	500	LYS

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Mol	Chain	Res	Type
1	B	503	ASN
1	B	506	THR
1	B	532	LEU
1	B	540	THR
1	B	544	LEU
1	B	554	PHE
1	B	556	GLU
1	B	557	LYS
1	B	571	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	543	GLN
1	B	543	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)

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 (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	578/585 (98%)	0.31	30 (5%)	27	34	32, 50, 89, 127	0
1	B	578/585 (98%)	0.29	35 (6%)	21	27	29, 49, 90, 121	0
All	All	1156/1170 (98%)	0.30	65 (5%)	24	30	29, 49, 90, 127	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	575	LEU	13.1
1	A	581	ALA	10.3
1	B	582	ALA	10.0
1	A	575	LEU	9.2
1	A	576	VAL	9.0
1	A	85	GLY	8.8
1	A	577	ALA	7.7
1	A	78	ALA	7.3
1	B	82	GLU	7.2
1	A	580	GLN	7.1
1	B	578	ALA	6.9
1	A	79	THR	6.7
1	A	83	THR	6.6
1	B	576	VAL	6.4
1	A	582	ALA	6.4
1	B	581	ALA	6.3
1	A	82	GLU	6.2
1	B	362	ALA	5.8
1	B	83	THR	5.6
1	B	579	SER	5.4
1	A	84	TYR	4.9
1	B	80	LEU	4.7
1	A	80	LEU	4.6
1	A	87	MET	4.6

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Mol	Chain	Res	Type	RSRZ
1	B	81	ARG	4.5
1	A	579	SER	3.6
1	A	574	LYS	3.6
1	A	539	ALA	3.5
1	B	79	THR	3.4
1	B	85	GLY	3.4
1	B	88	ALA	3.3
1	B	84	TYR	3.3
1	B	551	PHE	3.2
1	B	87	MET	3.2
1	A	362	ALA	3.1
1	B	365	ASP	3.1
1	B	574	LYS	3.0
1	A	572	GLY	2.9
1	A	81	ARG	2.9
1	B	364	ALA	2.9
1	B	580	GLN	2.8
1	B	367	HIS	2.8
1	A	578	ALA	2.8
1	B	77	VAL	2.7
1	B	86	GLU	2.6
1	A	503	ASN	2.6
1	A	538	LYS	2.6
1	B	503	ASN	2.5
1	B	283	LEU	2.5
1	B	535	HIS	2.5
1	B	539	ALA	2.5
1	B	502	PHE	2.4
1	A	551	PHE	2.4
1	A	86	GLU	2.4
1	B	370	TYR	2.4
1	A	547	VAL	2.3
1	B	544	LEU	2.3
1	A	519	LYS	2.3
1	A	560	LYS	2.2
1	B	519	LYS	2.2
1	B	570	GLU	2.2
1	A	573	LYS	2.1
1	B	78	ALA	2.1
1	B	577	ALA	2.0
1	A	535	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.