



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 04:40 am BST

PDB ID : 4EMX
Title : Crystal structure analysis of Human Serum Albumin in complex with chloride anions at cryogenic temperature
Authors : Botti, H.; Bonilla, L.; Trajtenberg, F.; Radi, R.; Buschiazzo, A.
Deposited on : 2012-04-12
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 ⓘ symbol.

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 : 2.11
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) : 2.11

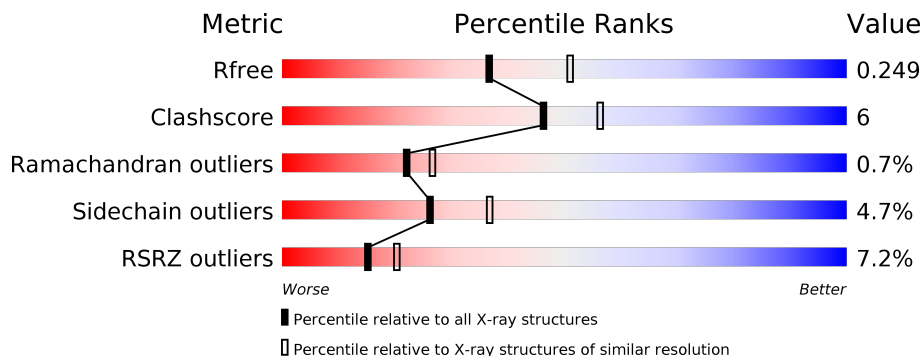
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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 on 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 $\leq 5\%$. 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	585	
1	B	585	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9377 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	581	4600	2903	784	872	41	0	5	0
1	B	579	4543	2867	772	863	41	0	4	0

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Cl	0	0
			3	3		
2	A	3	Total	Cl	0	0
			3	3		

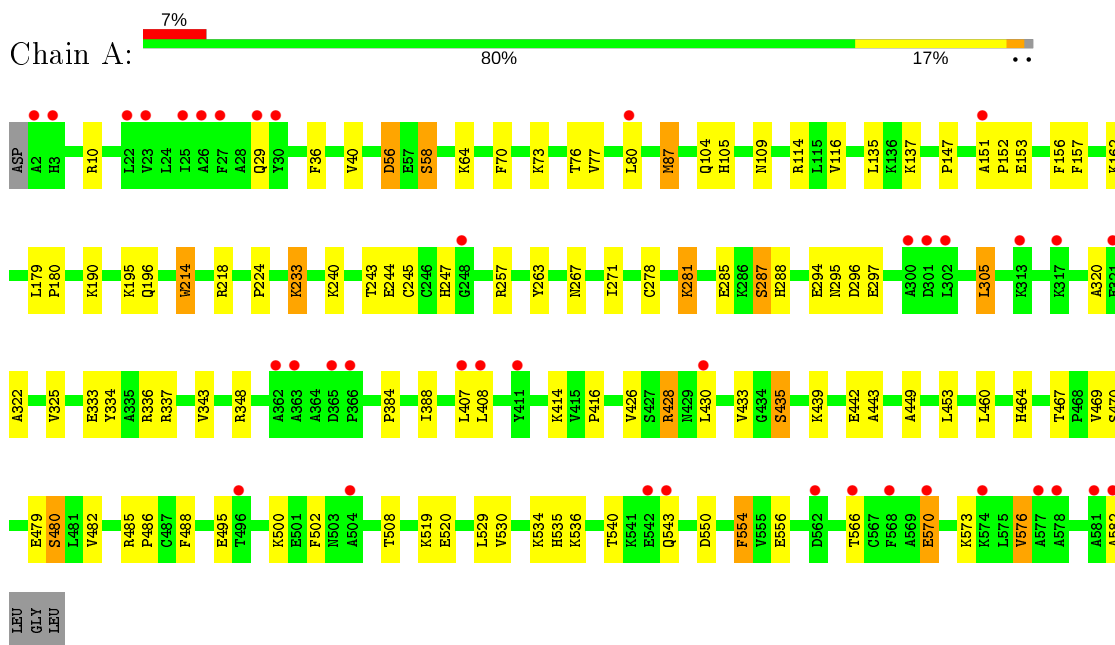
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	123	Total	O	0	0
			123	123		
3	B	105	Total	O	0	0
			105	105		

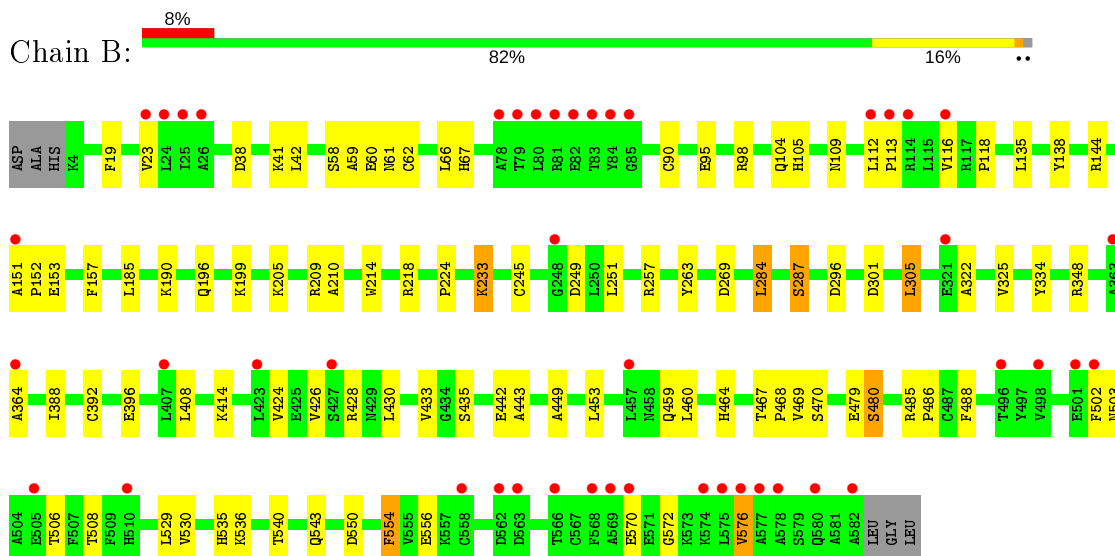
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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



- Molecule 1: Serum albumin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.88Å 59.23Å 95.61Å 75.17° 87.83° 74.20°	Depositor
Resolution (Å)	25.50 – 2.30 24.51 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.0 (25.50-2.30) 93.1 (24.51-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.31Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.187 , 0.238 0.192 , 0.249	Depositor DCC
R_{free} test set	1009 reflections (2.03%)	wwPDB-VP
Wilson B-factor (Å ²)	44.8	Xtrriage
Anisotropy	0.063	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 61.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9377	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section:
CL

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	0.75	1/4698 (0.0%)	0.85	6/6344 (0.1%)
1	B	0.71	1/4639 (0.0%)	0.83	3/6270 (0.0%)
All	All	0.73	2/9337 (0.0%)	0.84	9/12614 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	214	TRP	CD2-CE2	6.57	1.49	1.41
1	A	214	TRP	CD2-CE2	5.25	1.47	1.41

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	218	ARG	NE-CZ-NH1	8.02	124.31	120.30
1	B	218	ARG	NE-CZ-NH2	-7.65	116.47	120.30
1	A	428	ARG	NE-CZ-NH2	-6.58	117.01	120.30
1	A	348	ARG	NE-CZ-NH1	6.43	123.52	120.30
1	A	348	ARG	NE-CZ-NH2	-6.19	117.21	120.30
1	B	348	ARG	NE-CZ-NH2	-6.17	117.22	120.30
1	A	87	MET	CG-SD-CE	-5.88	90.79	100.20
1	A	10	ARG	NE-CZ-NH2	-5.72	117.44	120.30
1	A	56	ASP	CB-CG-OD1	5.27	123.05	118.30

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	4600	0	4487	66	0
1	B	4543	0	4408	56	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	123	0	0	5	0
3	B	105	0	0	5	0
All	All	9377	0	8895	113	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 6.

All (113) 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:500:LYS:NZ	3:A:779:HOH:O	2.04	0.88
1:A:464:HIS:HE1	1:A:470:SER:H	1.22	0.84
1:B:464:HIS:HE1	1:B:470:SER:H	1.27	0.81
1:A:320:ALA:HA	1:B:364:ALA:HB1	1.63	0.79
1:A:104:GLN:HG2	3:A:753:HOH:O	1.84	0.78
1:B:430:LEU:O	1:B:433:VAL:HG12	1.84	0.77
1:A:156:PHE:HE1	1:A:285:GLU:HG2	1.48	0.77
1:B:284:LEU:HD22	3:B:766:HOH:O	1.89	0.71
1:B:112:LEU:HD11	1:B:144:ARG:NH2	2.06	0.70
1:B:464:HIS:CE1	1:B:470:SER:H	2.08	0.70
1:A:305:LEU:HD22	1:A:334:TYR:HD2	1.56	0.69
1:A:320:ALA:HA	1:B:364:ALA:CB	2.23	0.69
1:A:156:PHE:CE1	1:A:285:GLU:HG2	2.27	0.68
1:A:64:LYS:HZ1	1:B:104:GLN:NE2	1.90	0.68
1:B:67:HIS:CE1	1:B:249:ASP:OD1	2.47	0.68
1:B:224:PRO:HD2	1:B:296:ASP:HB3	1.78	0.65
1:A:56:ASP:OD1	1:A:58:SER:HB2	1.97	0.64
1:A:464:HIS:CE1	1:A:470:SER:H	2.11	0.63
1:B:305:LEU:HD22	1:B:334:TYR:HD2	1.63	0.63
1:A:479:GLU:O	1:A:480:SER:HB2	1.99	0.63
1:A:64:LYS:NZ	1:B:104:GLN:NE2	2.46	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:ALA:HB3	1:B:152:PRO:CD	2.30	0.62
1:A:479:GLU:O	1:A:480:SER:CB	2.48	0.62
1:B:90:CYS:O	1:B:98:ARG:HG3	2.01	0.61
1:A:64:LYS:HZ1	1:B:104:GLN:HE22	1.49	0.60
1:A:414:LYS:HE3	1:A:488:PHE:O	2.02	0.59
1:A:243:THR:O	1:A:247[B]:HIS:HD2	1.87	0.57
1:B:479:GLU:O	1:B:480:SER:CB	2.52	0.57
1:A:29:GLN:HG2	1:A:147:PRO:HA	1.88	0.56
1:B:459:GLN:NE2	3:B:772:HOH:O	2.34	0.56
1:A:243:THR:O	1:A:247[B]:HIS:CD2	2.60	0.55
1:A:257:ARG:CZ	1:A:287:SER:HB3	2.36	0.55
1:A:464:HIS:CE1	1:A:469:VAL:H	2.24	0.55
1:B:257:ARG:CZ	1:B:287:SER:HB3	2.37	0.54
1:B:414:LYS:HE3	1:B:488:PHE:O	2.08	0.54
1:A:224:PRO:HD2	1:A:296:ASP:HB3	1.90	0.53
1:A:278:CYS:O	1:A:281[A]:LYS:HG3	2.08	0.53
1:B:67:HIS:HE1	1:B:249:ASP:OD1	1.90	0.52
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.90	0.52
1:A:502:PHE:HB2	1:A:535:HIS:CE1	2.44	0.52
1:B:153:GLU:HG3	1:B:157:PHE:HE1	1.74	0.52
1:A:430:LEU:O	1:A:433:VAL:HG12	2.10	0.52
1:B:424:VAL:O	1:B:428:ARG:HG3	2.10	0.51
1:A:408:LEU:HD22	1:A:529:LEU:HD23	1.92	0.50
1:A:573:LYS:HA	1:A:576:VAL:HG13	1.93	0.50
1:B:550:ASP:O	1:B:554:PHE:HB3	2.12	0.50
1:A:87:MET:CE	1:A:105:HIS:HB3	2.41	0.50
1:A:257:ARG:NH2	3:A:823:HOH:O	2.35	0.49
1:B:449:ALA:O	1:B:453:LEU:HG	2.13	0.49
1:A:153:GLU:OE2	1:A:288:HIS:ND1	2.39	0.48
1:A:322:ALA:HB1	1:A:325:VAL:HB	1.95	0.48
1:A:135:LEU:HD11	1:A:162:LYS:HB2	1.95	0.48
1:A:550:ASP:O	1:A:554:PHE:HB3	2.12	0.48
1:B:19:PHE:O	1:B:23:VAL:HG23	2.13	0.48
1:B:269[B]:ASP:OD1	1:B:269[B]:ASP:N	2.47	0.47
1:B:408:LEU:HD22	1:B:529:LEU:HD23	1.95	0.47
1:B:205:LYS:HD3	3:B:804:HOH:O	2.16	0.46
1:A:408:LEU:HD21	1:A:530:VAL:CG2	2.46	0.46
1:B:502:PHE:HB2	1:B:535:HIS:CE1	2.50	0.46
1:A:543:GLN:NE2	1:A:582:ALA:O	2.48	0.46
1:B:540:THR:O	1:B:543:GLN:N	2.49	0.46
1:A:320:ALA:CA	1:B:364:ALA:HB1	2.39	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:LEU:HD13	1:B:251:LEU:HD12	1.97	0.46
1:A:384:PRO:O	1:A:388:ILE:HG12	2.16	0.45
1:B:190:LYS:HB3	1:B:190:LYS:HE2	1.73	0.45
1:B:153:GLU:HG3	1:B:157:PHE:CE1	2.52	0.45
1:B:388:ILE:HD13	1:B:388:ILE:N	2.32	0.45
1:B:60:GLU:O	1:B:61[B]:ASN:HB2	2.16	0.45
1:B:392:CYS:O	1:B:396:GLU:HG2	2.17	0.45
1:B:151:ALA:HB3	1:B:152:PRO:HD3	1.98	0.45
1:A:190:LYS:HE2	1:A:190:LYS:HB3	1.69	0.45
1:B:572:GLY:O	1:B:576:VAL:HG23	2.17	0.45
1:B:116:VAL:O	1:B:118:PRO:HD3	2.18	0.44
1:A:233:LYS:HE3	1:A:263:TYR:CZ	2.52	0.44
1:A:240:LYS:O	1:A:244:GLU:HG3	2.16	0.44
1:A:428:ARG:HH12	1:A:519:LYS:NZ	2.16	0.44
1:A:196:GLN:HA	1:A:196:GLN:NE2	2.32	0.44
1:A:73:LYS:HA	1:A:73:LYS:HD3	1.89	0.44
1:B:209:ARG:HB3	3:B:734:HOH:O	2.17	0.44
1:A:179:LEU:N	1:A:180:PRO:CD	2.80	0.44
1:A:267:ASN:O	1:A:271:ILE:HG13	2.18	0.43
1:A:195:LYS:NZ	3:A:782:HOH:O	2.51	0.43
1:A:408:LEU:HD21	1:A:530:VAL:HG23	1.99	0.43
1:B:59:ALA:HB3	1:B:62:CYS:SG	2.58	0.43
1:A:214:TRP:CD1	1:A:343:VAL:HG11	2.53	0.43
1:A:76:THR:HA	1:B:95:GLU:O	2.19	0.43
1:A:218:ARG:NH2	1:A:295:ASN:OD1	2.49	0.43
1:A:36:PHE:O	1:A:40:VAL:HG23	2.18	0.43
1:B:210:ALA:N	3:B:734:HOH:O	2.50	0.43
1:B:196:GLN:O	1:B:199:LYS:N	2.51	0.43
1:B:322:ALA:HB1	1:B:325:VAL:HB	2.00	0.43
1:A:320:ALA:CA	1:B:364:ALA:CB	2.93	0.43
1:B:464:HIS:CE1	1:B:469:VAL:H	2.36	0.43
1:B:233:LYS:HE3	1:B:263:TYR:CE2	2.55	0.42
1:A:426:VAL:HG21	1:A:460:LEU:HB2	2.00	0.42
1:A:196:GLN:NE2	3:A:773:HOH:O	2.52	0.42
1:A:153:GLU:O	1:A:157:PHE:HD1	2.02	0.42
1:A:114:ARG:NH1	1:A:520:GLU:HG2	2.34	0.42
1:B:408:LEU:HD21	1:B:530:VAL:HG23	2.01	0.42
1:A:333:GLU:HB3	1:A:337:ARG:NH1	2.34	0.42
1:A:416:PRO:O	1:A:534:LYS:HE2	2.19	0.42
1:A:70:PHE:N	1:A:70:PHE:CD1	2.87	0.42
1:A:407:LEU:HD13	1:A:430:LEU:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:SER:O	1:A:439:LYS:HE3	2.20	0.41
1:B:135:LEU:O	1:B:138:TYR:HB3	2.21	0.41
1:B:426:VAL:HG21	1:B:460:LEU:HB2	2.03	0.41
1:B:503:ASN:HB3	1:B:506:THR:HB	2.03	0.41
1:A:87:MET:HE2	1:A:105:HIS:HB3	2.03	0.40
1:A:151:ALA:N	1:A:152:PRO:HD2	2.36	0.40
1:B:38:ASP:O	1:B:42:LEU:HG	2.21	0.40
1:A:449:ALA:O	1:A:453:LEU:HG	2.22	0.40
1:A:320:ALA:C	1:B:364:ALA:CB	2.89	0.40
1:B:485:ARG:HB3	1:B:486:PRO:HD3	2.03	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	584/585 (100%)	562 (96%)	18 (3%)	4 (1%)	22	26
1	B	581/585 (99%)	555 (96%)	22 (4%)	4 (1%)	22	26
All	All	1165/1170 (100%)	1117 (96%)	40 (3%)	8 (1%)	22	26

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	443	ALA
1	A	480	SER
1	B	443	ALA
1	B	480	SER
1	A	442	GLU
1	B	442	GLU
1	B	576	VAL
1	A	570	GLU

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	497/511 (97%)	470 (95%)	27 (5%)	22	30
1	B	488/511 (96%)	468 (96%)	20 (4%)	30	43
All	All	985/1022 (96%)	938 (95%)	47 (5%)	26	36

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	SER
1	A	77	VAL
1	A	80	LEU
1	A	109	ASN
1	A	116	VAL
1	A	137	LYS
1	A	233	LYS
1	A	245	CYS
1	A	281[A]	LYS
1	A	281[B]	LYS
1	A	287	SER
1	A	294	GLU
1	A	297	GLU
1	A	305	LEU
1	A	336	ARG
1	A	435	SER
1	A	467	THR
1	A	482	VAL
1	A	495	GLU
1	A	508	THR
1	A	536	LYS
1	A	540	THR
1	A	554	PHE
1	A	556	GLU
1	A	566	THR
1	A	570	GLU
1	A	576	VAL

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Mol	Chain	Res	Type
1	B	41	LYS
1	B	58	SER
1	B	105	HIS
1	B	109	ASN
1	B	113	PRO
1	B	185	LEU
1	B	233	LYS
1	B	245	CYS
1	B	284	LEU
1	B	287	SER
1	B	301	ASP
1	B	305	LEU
1	B	435	SER
1	B	467	THR
1	B	468	PRO
1	B	508	THR
1	B	536	LYS
1	B	554	PHE
1	B	556	GLU
1	B	570	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	ASN
1	A	196	GLN
1	A	464	HIS
1	A	483	ASN
1	B	33	GLN
1	B	67	HIS
1	B	104	GLN
1	B	109	ASN
1	B	196	GLN
1	B	247	HIS
1	B	464	HIS
1	B	483	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 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

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	581/585 (99%)	0.28	39 (6%) 17 23	43, 69, 129, 178	0
1	B	579/585 (98%)	0.28	45 (7%) 13 17	48, 72, 129, 173	0
All	All	1160/1170 (99%)	0.28	84 (7%) 15 20	43, 71, 129, 178	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	78	ALA	7.7
1	B	84	TYR	5.8
1	B	577	ALA	5.8
1	B	83	THR	5.6
1	B	575	LEU	5.2
1	B	79	THR	5.1
1	A	317	LYS	5.0
1	A	2	ALA	4.6
1	B	80	LEU	4.3
1	B	582	ALA	4.2
1	B	25	ILE	4.1
1	A	582	ALA	4.0
1	A	3	HIS	4.0
1	A	562	ASP	3.9
1	B	363	ALA	3.7
1	A	25	ILE	3.6
1	B	578	ALA	3.6
1	B	85	GLY	3.5
1	A	581	ALA	3.5
1	B	568	PHE	3.5
1	B	82	GLU	3.4
1	A	570	GLU	3.4
1	A	321	GLU	3.3
1	A	578	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	364	ALA	3.2
1	A	362	ALA	3.2
1	A	577	ALA	3.2
1	A	504	ALA	3.1
1	B	574	LYS	3.1
1	B	563	ASP	3.0
1	B	569	ALA	3.0
1	A	542	GLU	3.0
1	B	113	PRO	3.0
1	A	26	ALA	2.9
1	B	501	GLU	2.9
1	B	26	ALA	2.9
1	B	558	CYS	2.9
1	A	543	GLN	2.8
1	B	510	HIS	2.8
1	B	81	ARG	2.7
1	B	496	THR	2.6
1	A	22	LEU	2.6
1	B	423	LEU	2.6
1	A	363	ALA	2.6
1	A	408	LEU	2.5
1	B	580	GLN	2.5
1	A	300	ALA	2.5
1	A	301	ASP	2.5
1	A	430	LEU	2.5
1	A	566	THR	2.4
1	B	562	ASP	2.4
1	A	407	LEU	2.4
1	B	576	VAL	2.4
1	A	411	TYR	2.4
1	B	112	LEU	2.4
1	B	570	GLU	2.4
1	A	151	ALA	2.4
1	A	365	ASP	2.3
1	B	502	PHE	2.3
1	A	313	LYS	2.3
1	A	568	PHE	2.3
1	B	407	LEU	2.3
1	B	24	LEU	2.2
1	B	114	ARG	2.2
1	A	574	LYS	2.2
1	B	248	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	151	ALA	2.2
1	A	30	TYR	2.2
1	A	248	GLY	2.2
1	B	498	VAL	2.2
1	A	496	THR	2.2
1	A	23	VAL	2.1
1	A	302	LEU	2.1
1	A	366	PRO	2.1
1	B	427	SER	2.1
1	B	505	GLU	2.1
1	A	27	PHE	2.1
1	B	566	THR	2.1
1	A	29	GLN	2.0
1	B	23	VAL	2.0
1	B	116	VAL	2.0
1	B	321	GLU	2.0
1	A	80	LEU	2.0
1	B	457	LEU	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](#)

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
2	CL	A	601	1/1	0.93	0.06	98,98,98,98	0
2	CL	B	602	1/1	0.93	0.15	81,81,81,81	0
2	CL	B	601	1/1	0.94	0.33	80,80,80,80	0
2	CL	A	603	1/1	0.98	0.43	92,92,92,92	0
2	CL	A	602	1/1	0.99	0.11	67,67,67,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	CL	B	603	1/1	0.99	0.20	78,78,78,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.