



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:54 PM GMT

PDB ID : 3S7I
Title : Crystal structure of Ara h 1
Authors : Chruszcz, M.; Maleki, S.J.; Solberg, R.; Minor, W.
Deposited on : 2011-05-26
Resolution : 2.35 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

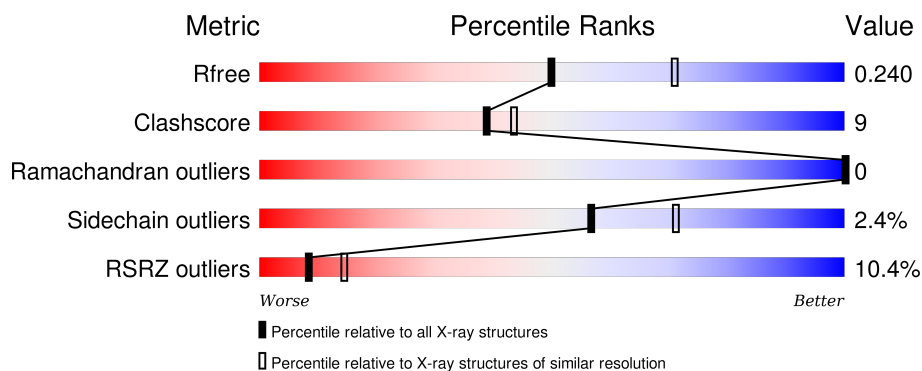
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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	418	<div> <div>17%</div> <div>72%</div> <div>16%</div> <div>12%</div> </div>
1	B	418	<div> <div>69%</div> <div>17%</div> <div>13%</div> </div>

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
2	CL	A	1	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5731 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 Allergen Ara h 1, clone P41B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	S	0	1	0
			2898	1813	526	552	7			
1	B	363	Total	C	N	O	S	0	0	0
			2698	1696	480	515	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	169	MET	-	EXPRESSION TAG	UNP P43238
B	169	MET	-	EXPRESSION TAG	UNP P43238

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		

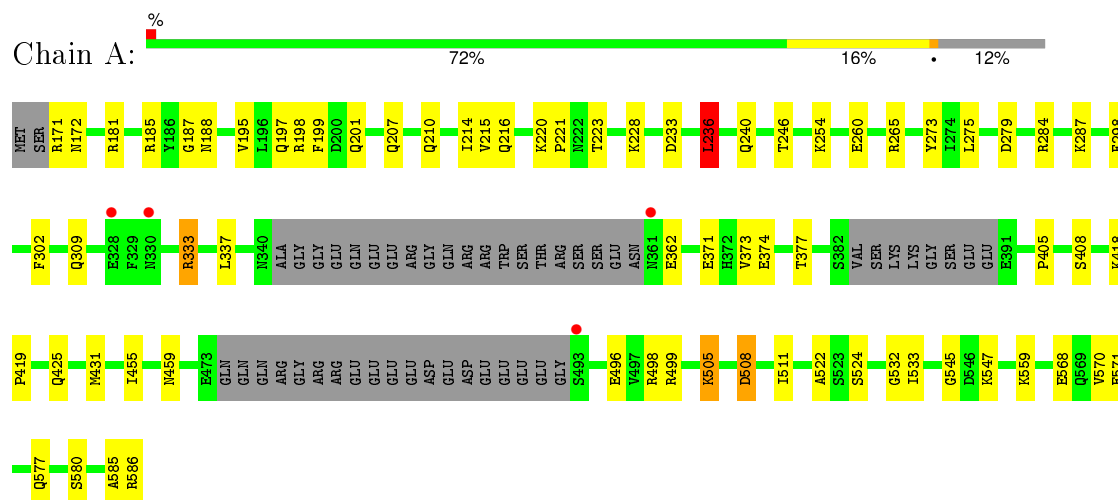
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	114	Total	O	0	0
			114	114		
3	B	19	Total	O	0	0
			19	19		

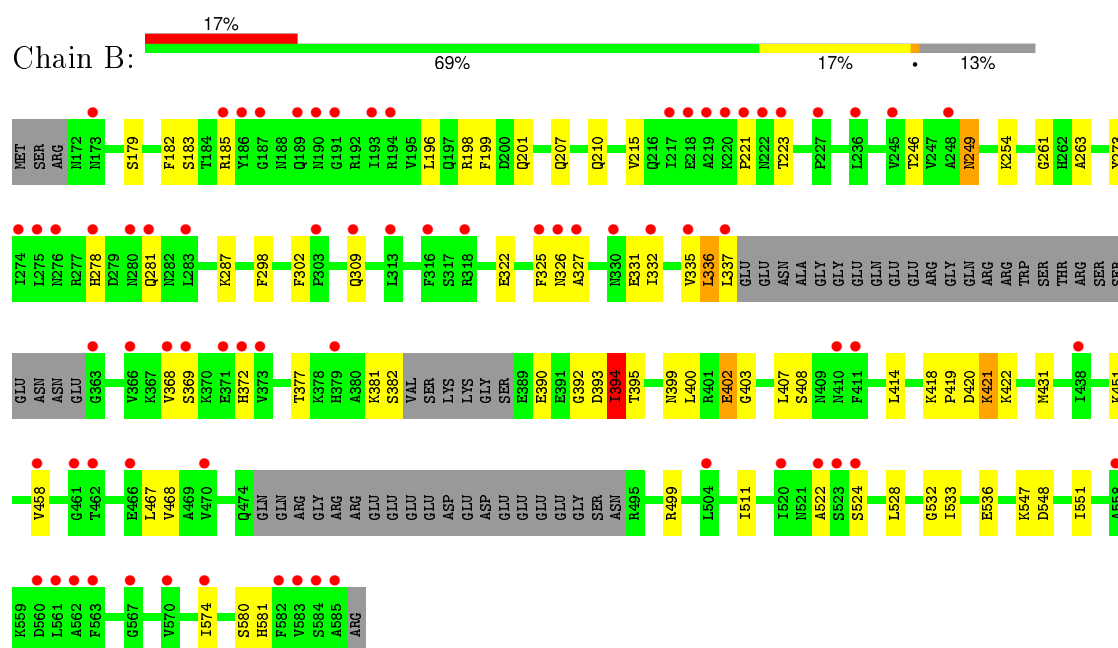
3 Residue-property plots

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: Allergen Ara h 1, clone P41B



• Molecule 1: Allergen Ara h 1, clone P41B



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	92.89Å 92.89Å 231.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.13 – 2.35 40.13 – 2.35	Depositor EDS
% Data completeness (in resolution range)	86.5 (40.13-2.35) 99.9 (40.13-2.35)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.14 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.203 , 0.244 0.219 , 0.240	Depositor DCC
R_{free} test set	1566 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	44.1	Xtriage
Anisotropy	0.889	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.8	EDS
Estimated twinning fraction	0.948 for H, K, L 0.052 for K, H, -L 0.008 for -1/3*h+1/3*k+1/3*l,-k,8/3*h+4/ 3*k+1/3*l 0.019 for -2/3*h-1/3*k-1/3*l,-1/3*h-2/3*k+ 1/3*l,-4/3*h+4/3*k+1/3*l 0.003 for -h,1/3*h-1/3*k-1/3*l,-4/3*h-8/3*k +1/3*l 0.007 for -h,2/3*h+1/3*k+1/3*l,4/3*h+8/3 *k-1/3*l 0.006 for 1/3*h+2/3*k-1/3*l,-k,-8/3*h-4/3* k-1/3*l 0.016 for -1/3*h-2/3*k+1/3*l,-2/3*h-1/3*k- 1/3*l,4/3*h-4/3*k-1/3*l 0.037 for -h-k,k,-l	Xtriage
Reported twinning fraction	0.948 for H, K, L 0.052 for K, H, -L	Depositor
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 31041 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5731	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% of the height of the origin peak. No significant pseudotranslation is detected.*

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	1.06	2/2947 (0.1%)	0.92	2/3980 (0.1%)
1	B	0.60	0/2746	0.72	1/3734 (0.0%)
All	All	0.87	2/5693 (0.0%)	0.83	3/7714 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	254	LYS	CE-NZ	5.75	1.63	1.49
1	A	260	GLU	CB-CG	-5.55	1.41	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	508	ASP	CB-CG-OD2	5.48	123.23	118.30
1	B	394	ILE	N-CA-C	-5.30	96.69	111.00
1	A	236	LEU	CA-CB-CG	5.17	127.19	115.30

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	2898	0	2816	41	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2698	0	2506	58	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	114	0	0	5	0
3	B	19	0	0	0	0
All	All	5731	0	5322	98	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 9.

All (98) 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:336:LEU:O	1:B:337:LEU:HB2	1.49	1.11
1:B:421:LYS:H	1:B:421:LYS:HE3	1.37	0.88
1:B:392:GLY:O	1:B:393:ASP:HB2	1.76	0.86
1:B:421:LYS:H	1:B:421:LYS:CE	1.91	0.83
1:B:207:GLN:O	1:B:210:GLN:HG2	1.84	0.78
1:B:336:LEU:O	1:B:337:LEU:CB	2.30	0.76
1:B:368:VAL:HG12	1:B:372:HIS:HB2	1.71	0.71
1:B:399:ASN:O	1:B:402:GLU:HB3	1.90	0.71
1:A:333:ARG:HA	1:A:337:LEU:HB2	1.76	0.66
1:B:369:SER:HB2	1:B:372:HIS:ND1	2.11	0.65
1:A:199:PHE:CD1	1:A:511:ILE:HD11	2.30	0.65
1:B:263:ALA:HB2	1:B:400:LEU:HD21	1.79	0.64
1:B:431:MET:HG3	1:B:532:GLY:O	1.98	0.64
1:A:505:LYS:NZ	1:A:508:ASP:OD2	2.31	0.64
1:A:585:ALA:O	1:A:586:ARG:CB	2.47	0.63
1:A:197:GLN:HB3	3:A:51:HOH:O	1.97	0.63
1:B:199:PHE:CD1	1:B:511:ILE:HD11	2.34	0.63
1:B:421:LYS:N	1:B:421:LYS:HE3	2.12	0.62
1:B:399:ASN:HB3	1:B:402:GLU:HB2	1.81	0.61
1:A:498:ARG:NH1	1:A:499:ARG:O	2.33	0.61
1:A:431:MET:HG3	1:A:532:GLY:O	2.03	0.59
1:A:207:GLN:O	1:A:210:GLN:HG2	2.03	0.59
1:A:568:GLU:HG3	3:A:76:HOH:O	2.05	0.57
1:B:467:LEU:HD23	1:B:468:VAL:N	2.18	0.57
1:B:325:PHE:O	1:B:326:ASN:HB3	2.04	0.57
1:A:533:ILE:HD12	1:A:533:ILE:N	2.20	0.57
1:B:377:THR:O	1:B:377:THR:HG22	2.05	0.56
1:B:421:LYS:N	1:B:421:LYS:CE	2.65	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:261:GLY:O	1:B:400:LEU:HG	2.07	0.55
1:A:185:ARG:NH1	1:A:298:PHE:O	2.36	0.55
1:A:188:ASN:ND2	1:A:362:GLU:O	2.38	0.54
1:A:547:LYS:HB2	1:A:580:SER:OG	2.08	0.53
1:A:221:PRO:O	1:A:223:THR:HG23	2.09	0.53
1:A:171:ARG:N	1:B:183:SER:HG	2.07	0.53
1:B:421:LYS:H	1:B:421:LYS:HE2	1.73	0.52
1:A:187:GLY:O	1:A:188:ASN:HB3	2.10	0.52
1:B:331:GLU:O	1:B:335:VAL:HG23	2.09	0.52
1:A:198:ARG:HB2	1:A:201:GLN:HG3	1.90	0.52
1:B:399:ASN:HB3	1:B:402:GLU:CB	2.40	0.52
1:A:199:PHE:CE1	1:A:511:ILE:HD11	2.45	0.51
1:A:284:ARG:NH2	1:A:459:ASN:O	2.42	0.51
1:B:278:HIS:CD2	1:B:281:GLN:HB3	2.45	0.50
1:B:185:ARG:NH1	1:B:298:PHE:O	2.40	0.50
1:A:559:LYS:HE3	1:A:571:GLU:HG2	1.94	0.50
1:B:522:ALA:HB1	1:B:524:SER:O	2.12	0.50
1:A:236:LEU:C	1:A:236:LEU:HD23	2.32	0.50
1:B:548:ASP:OD1	1:B:581:HIS:NE2	2.45	0.49
1:B:390:GLU:HB2	1:B:395:THR:HA	1.94	0.49
1:B:246:THR:HG21	1:B:273:TYR:CZ	2.47	0.49
1:A:215:VAL:HB	1:A:287:LYS:HB2	1.94	0.49
1:B:402:GLU:HG3	1:B:403:GLY:N	2.23	0.48
1:B:532:GLY:C	1:B:533:ILE:HD12	2.33	0.48
1:B:533:ILE:HD12	1:B:533:ILE:N	2.27	0.48
1:A:559:LYS:HD2	1:A:570:VAL:HG11	1.96	0.48
1:A:302:PHE:CD1	1:A:309:GLN:HG3	2.49	0.47
1:B:302:PHE:CD1	1:B:309:GLN:HG3	2.51	0.46
1:A:545:GLY:HA2	1:A:577:GLN:HG3	1.96	0.46
1:B:368:VAL:CG1	1:B:372:HIS:CB	2.94	0.46
1:A:220:LYS:HB3	1:A:221:PRO:HD2	1.97	0.45
1:A:240:GLN:HB3	1:A:284:ARG:HB2	1.98	0.45
1:B:368:VAL:CG1	1:B:372:HIS:HB2	2.42	0.45
1:A:371:GLU:O	1:A:374:GLU:HB2	2.16	0.45
1:A:185:ARG:HH21	1:A:195:VAL:HG21	1.81	0.45
1:A:418:LYS:HB3	1:A:419:PRO:HD2	1.98	0.45
1:B:400:LEU:N	1:B:400:LEU:HD23	2.32	0.45
1:B:369:SER:CB	1:B:372:HIS:ND1	2.79	0.44
1:B:182:PHE:CE1	1:B:196:LEU:HD13	2.52	0.44
1:B:421:LYS:HE3	1:B:421:LYS:HB2	1.53	0.44
1:B:467:LEU:C	1:B:467:LEU:HD23	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:ARG:HD3	1:A:425:GLN:OE1	2.18	0.44
1:B:381:LYS:O	1:B:382:SER:CB	2.65	0.43
1:B:215:VAL:HB	1:B:287:LYS:HB2	2.00	0.43
1:B:547:LYS:N	1:B:580:SER:OG	2.51	0.43
1:A:181:ARG:O	3:A:51:HOH:O	2.21	0.43
1:A:522:ALA:HB1	1:A:524:SER:O	2.19	0.43
1:A:233:ASP:CG	1:A:265:ARG:HH11	2.21	0.42
1:B:249:ASN:OD1	1:B:394:ILE:HD11	2.19	0.42
1:A:405:PRO:HG2	1:A:408:SER:HB2	2.00	0.42
1:B:337:LEU:HD23	1:B:337:LEU:HA	1.82	0.42
1:B:451:LYS:HD3	1:B:536:GLU:O	2.19	0.42
1:B:551:ILE:HG22	1:B:574:ILE:HD11	2.02	0.42
1:B:418:LYS:HB3	1:B:419:PRO:HD2	2.02	0.41
1:B:278:HIS:HD2	1:B:281:GLN:CB	2.32	0.41
1:B:221:PRO:O	1:B:223:THR:HG23	2.20	0.41
1:B:414:LEU:C	1:B:414:LEU:HD23	2.40	0.41
1:B:198:ARG:HB2	1:B:201:GLN:HG3	2.02	0.41
1:A:214:ILE:HD11	1:A:455:ILE:HG21	2.02	0.41
1:B:254:LYS:NZ	1:B:390:GLU:HA	2.35	0.41
1:B:332:ILE:O	1:B:336:LEU:O	2.39	0.41
1:B:322:GLU:HG2	1:B:327:ALA:O	2.21	0.41
1:B:420:ASP:HB3	1:B:422:LYS:H	1.86	0.41
1:A:216:GLN:HB3	3:A:139:HOH:O	2.21	0.41
1:A:228:LYS:HE3	3:A:59:HOH:O	2.21	0.41
1:B:407:LEU:HD22	1:B:581:HIS:CE1	2.56	0.40
1:A:246:THR:HG21	1:A:273:TYR:CZ	2.56	0.40
1:B:458:VAL:HG22	1:B:528:LEU:CD2	2.52	0.40
1:A:171:ARG:HB2	1:A:172:ASN:H	1.70	0.40
1:A:275:LEU:HD21	1:A:373:VAL:HG13	2.04	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	362/418 (87%)	355 (98%)	7 (2%)	0	100	100
1	B	355/418 (85%)	346 (98%)	9 (2%)	0	100	100
All	All	717/836 (86%)	701 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	311/365 (85%)	305 (98%)	6 (2%)	65	79
1	B	271/365 (74%)	263 (97%)	8 (3%)	48	62
All	All	582/730 (80%)	568 (98%)	14 (2%)	57	71

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	236	LEU
1	A	279	ASP
1	A	333	ARG
1	A	377	THR
1	A	496	GLU
1	A	505	LYS
1	B	179	SER
1	B	249	ASN
1	B	336	LEU
1	B	394	ILE
1	B	402	GLU
1	B	408	SER
1	B	421	LYS
1	B	499	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	396	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 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

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	369/418 (88%)	0.14	4 (1%) 82 90	13, 31, 54, 67	2 (0%)
1	B	363/418 (86%)	1.14	72 (19%) 1 2	40, 79, 117, 140	1 (0%)
All	All	732/836 (87%)	0.63	76 (10%) 8 14	13, 48, 112, 140	3 (0%)

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	316	PHE	6.5
1	B	461	GLY	5.9
1	B	275	LEU	5.6
1	B	219	ALA	5.6
1	B	373	VAL	5.2
1	B	335	VAL	5.1
1	B	327	ALA	5.0
1	B	524	SER	5.0
1	B	283	LEU	5.0
1	B	186	TYR	5.0
1	B	190	ASN	4.9
1	B	371	GLU	4.9
1	B	411	PHE	4.8
1	B	368	VAL	4.7
1	B	189	GLN	4.6
1	B	570	VAL	4.6
1	B	332	ILE	4.4
1	B	520	ILE	4.1
1	B	563	PHE	4.0
1	B	372	HIS	4.0
1	B	561	LEU	3.9
1	B	280	ASN	3.8
1	B	222	ASN	3.7
1	B	366	VAL	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	218	GLU	3.6
1	B	562	ALA	3.6
1	B	325	PHE	3.6
1	B	193	ILE	3.5
1	B	281	GLN	3.4
1	B	318	ARG	3.4
1	B	337	LEU	3.2
1	B	584	SER	3.1
1	B	236	LEU	3.1
1	B	369	SER	3.0
1	A	361	ASN	3.0
1	B	220	LYS	2.9
1	A	328	GLU	2.9
1	B	567	GLY	2.9
1	B	223	THR	2.9
1	B	410	ASN	2.8
1	B	574	ILE	2.8
1	B	470	VAL	2.8
1	B	363	GLY	2.8
1	B	245	VAL	2.7
1	B	458	VAL	2.7
1	A	330	ASN	2.7
1	B	313	LEU	2.7
1	B	221	PRO	2.7
1	B	438	ILE	2.6
1	B	278	HIS	2.6
1	B	326	ASN	2.6
1	B	583	VAL	2.6
1	B	173	ASN	2.6
1	B	309	GLN	2.5
1	B	504	LEU	2.5
1	B	558	ALA	2.5
1	B	462	THR	2.5
1	B	523	SER	2.5
1	B	303	PRO	2.5
1	B	379	HIS	2.4
1	B	522	ALA	2.4
1	B	274	ILE	2.4
1	B	248	ALA	2.4
1	B	187	GLY	2.4
1	B	185	ARG	2.4
1	A	493	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	191	GLY	2.3
1	B	276	ASN	2.3
1	B	582	PHE	2.2
1	B	585	ALA	2.2
1	B	560	ASP	2.1
1	B	227	PRO	2.1
1	B	217	ILE	2.1
1	B	466	GLU	2.0
1	B	330	ASN	2.0
1	B	194	ARG	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. 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	CL	A	1	1/1	0.81	0.35	17.58	56,56,56,56	0
2	CL	B	1	1/1	0.63	0.30	-	80,80,80,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.