



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 08:45 am GMT

PDB ID : 3A0R
Title : Crystal structure of histidine kinase ThkA (TM1359) in complex with response regulator protein TrrA (TM1360)
Authors : Yamada, S.; Sugimoto, H.; Kobayashi, M.; Ohno, A.; Nakamura, H.; Shiro, Y.
Deposited on : 2009-03-24
Resolution : 3.80 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

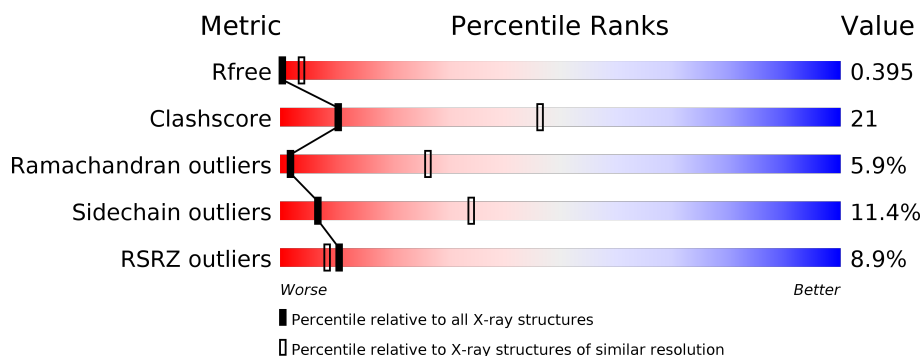
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 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}	100719	1019 (4.08-3.52)
Clashscore	112137	1030 (4.04-3.56)
Ramachandran outliers	110173	1011 (4.06-3.54)
Sidechain outliers	110143	1005 (4.06-3.54)
RSRZ outliers	101464	1032 (4.08-3.52)

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	349	
2	B	116	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3709 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 Sensor protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	0	0
			2768	1772	467	522	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	407	MET	-	EXPRESSION TAG	UNP Q9X180

- Molecule 2 is a protein called Response regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	116	Total	C	N	O	Se	0	0	0
			940	600	150	187	3			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	89	MSE	LEU	ENGINEERED	UNP Q9X181

- Molecule 3 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Hg	0	0
			1	1		

4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	110.59Å 110.59Å 352.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.80 48.98 – 3.80	Depositor EDS
% Data completeness (in resolution range)	98.2 (20.00-3.80) 98.1 (48.98-3.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.82 (at 3.77Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.350 , 0.371 0.353 , 0.395	Depositor DCC
R_{free} test set	546 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	152.5	Xtriage
Anisotropy	0.673	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 447.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3709	wwPDB-VP
Average B, all atoms (Å ²)	164.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.11% 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: HG

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	0.48	1/2814 (0.0%)	0.64	1/3780 (0.0%)
2	B	0.35	0/951	0.48	0/1269
All	All	0.45	1/3765 (0.0%)	0.61	1/5049 (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
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	721	CYS	CB-SG	5.78	1.92	1.82

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	658	LEU	CA-CB-CG	5.67	128.35	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	657	ASN	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	2768	0	2811	142	0
2	B	940	0	951	15	0
3	A	1	0	0	1	0
All	All	3709	0	3762	155	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 21.

All (155) 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:721:CYS:SG	3:A:800:HG:HG	1.47	1.33
1:A:653:GLN:HA	1:A:656:ILE:HD12	1.21	1.13
1:A:718:LEU:HD12	1:A:721:CYS:HB2	1.28	1.13
1:A:606:PHE:HD2	1:A:646:ALA:HA	1.36	0.90
1:A:659:VAL:O	1:A:663:ILE:HG13	1.72	0.88
1:A:633:PHE:CZ	1:A:674:ILE:HD11	2.08	0.87
1:A:417:GLU:HG3	1:A:419:ILE:H	1.39	0.85
1:A:606:PHE:CD2	1:A:646:ALA:HA	2.13	0.84
1:A:429:THR:O	1:A:436:ILE:HA	1.77	0.82
1:A:721:CYS:O	1:A:733:ILE:HD13	1.79	0.82
1:A:592:LYS:HA	1:A:595:LEU:HD12	1.59	0.82
1:A:653:GLN:O	1:A:656:ILE:HB	1.79	0.81
1:A:685:VAL:HB	1:A:746:PHE:CZ	2.18	0.79
2:B:52:ASP:HA	2:B:80:THR:HA	1.67	0.76
1:A:485:ASN:O	1:A:486:PHE:HB3	1.87	0.74
1:A:612:ASN:HA	1:A:615:ILE:HB	1.70	0.73
1:A:425:THR:O	1:A:442:LYS:HB3	1.89	0.73
1:A:718:LEU:CD1	1:A:721:CYS:HB2	2.12	0.72
1:A:633:PHE:HZ	1:A:674:ILE:HD11	1.53	0.71
1:A:550:ARG:HA	1:A:553:ILE:HB	1.73	0.69
2:B:5:LEU:HD11	2:B:40:LYS:HD3	1.74	0.69
1:A:591:VAL:O	1:A:595:LEU:HG	1.94	0.67
1:A:651:ILE:HG12	1:A:729:HIS:CE1	2.30	0.67
1:A:629:MET:O	1:A:631:ILE:HG13	1.95	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:655:LEU:O	1:A:656:ILE:C	2.33	0.65
1:A:653:GLN:CA	1:A:656:ILE:HD12	2.14	0.65
1:A:438:GLU:HA	1:A:451:LYS:HE2	1.79	0.64
1:A:546:ALA:HA	1:A:549:ILE:CD1	2.27	0.64
1:A:658:LEU:HD22	1:A:744:PHE:HB3	1.79	0.63
1:A:432:LYS:HA	1:A:476:PHE:CZ	2.33	0.63
1:A:647:ASP:O	1:A:651:ILE:HG13	1.98	0.63
1:A:546:ALA:HA	1:A:549:ILE:HD12	1.79	0.63
1:A:426:ALA:HA	1:A:440:ASN:HD22	1.62	0.63
1:A:633:PHE:CE1	1:A:674:ILE:HD11	2.35	0.62
1:A:485:ASN:HB3	1:A:604:LEU:HB2	1.83	0.61
1:A:549:ILE:HG22	1:A:553:ILE:HD12	1.81	0.61
1:A:417:GLU:HB3	1:A:420:LEU:HD23	1.83	0.61
1:A:485:ASN:O	1:A:486:PHE:CB	2.47	0.60
1:A:720:ILE:O	1:A:724:ILE:HG12	2.02	0.60
2:B:18:LYS:HG3	2:B:28:ILE:HB	1.83	0.60
1:A:644:VAL:HG12	1:A:750:LYS:HA	1.83	0.60
1:A:605:GLU:O	1:A:647:ASP:HA	2.01	0.59
1:A:674:ILE:HG23	1:A:687:VAL:HG22	1.83	0.59
1:A:542:THR:O	1:A:546:ALA:HB2	2.02	0.59
2:B:71:LYS:HB3	2:B:74:ALA:HB2	1.85	0.59
1:A:439:TRP:HB3	1:A:451:LYS:HG3	1.83	0.59
1:A:559:PHE:HA	1:A:562:ARG:HB2	1.84	0.58
1:A:734:TRP:HE3	1:A:734:TRP:H	1.52	0.58
1:A:658:LEU:HD22	1:A:744:PHE:CB	2.33	0.58
1:A:496:ILE:HG23	1:A:516:ILE:HG12	1.86	0.58
1:A:452:GLU:HA	1:A:455:LEU:HD12	1.86	0.58
1:A:700:GLU:HA	1:A:703:PHE:HD2	1.68	0.58
1:A:486:PHE:CG	1:A:603:VAL:HA	2.38	0.57
1:A:520:THR:C	1:A:522:LEU:H	2.08	0.57
1:A:646:ALA:CB	1:A:651:ILE:HD11	2.35	0.56
2:B:83:SER:HA	2:B:97:TYR:OH	2.06	0.56
1:A:703:PHE:HA	1:A:718:LEU:HD23	1.87	0.56
1:A:425:THR:O	1:A:440:ASN:ND2	2.39	0.55
1:A:552:PRO:HB2	1:A:584:LEU:HD22	1.88	0.55
1:A:420:LEU:O	1:A:422:SER:N	2.41	0.54
1:A:650:ARG:C	1:A:652:LYS:N	2.60	0.54
1:A:549:ILE:HG22	1:A:553:ILE:CD1	2.37	0.53
2:B:68:ARG:HH21	2:B:94:ALA:HA	1.74	0.53
1:A:700:GLU:O	1:A:703:PHE:HB2	2.09	0.53
1:A:497:ARG:HB3	1:A:515:THR:OG1	2.08	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:14:ARG:HE	2:B:30:THR:HB	1.74	0.53
1:A:502:ARG:HG2	1:A:508:LEU:O	2.09	0.53
1:A:429:THR:HB	1:A:438:GLU:OE1	2.08	0.53
1:A:615:ILE:HG22	1:A:616:ARG:N	2.24	0.52
1:A:725:ILE:HG21	1:A:732:LYS:O	2.10	0.52
1:A:592:LYS:HA	1:A:595:LEU:CD1	2.35	0.52
1:A:650:ARG:O	1:A:653:GLN:HG3	2.08	0.52
1:A:660:GLN:O	1:A:663:ILE:HB	2.10	0.52
1:A:486:PHE:CD1	1:A:603:VAL:HA	2.45	0.51
1:A:530:LYS:O	1:A:534:ARG:HD3	2.11	0.51
1:A:648:ARG:O	1:A:650:ARG:N	2.44	0.51
1:A:552:PRO:HG2	1:A:584:LEU:HD13	1.93	0.51
1:A:503:ASN:H	1:A:510:GLU:HG3	1.76	0.50
1:A:610:ASN:HA	1:A:642:LEU:O	2.12	0.50
1:A:720:ILE:H	1:A:720:ILE:HD12	1.77	0.50
2:B:47:ASP:O	2:B:75:LYS:HG3	2.11	0.50
1:A:650:ARG:C	1:A:652:LYS:H	2.14	0.49
1:A:542:THR:O	1:A:546:ALA:CB	2.60	0.49
1:A:590:ILE:HG22	1:A:590:ILE:O	2.12	0.49
1:A:531:ARG:HA	1:A:534:ARG:HB2	1.95	0.49
1:A:661:ASN:C	1:A:663:ILE:H	2.14	0.49
1:A:585:SER:O	1:A:588:GLU:HB2	2.13	0.49
1:A:492:ARG:HG2	1:A:520:THR:HG21	1.94	0.48
2:B:7:VAL:HG22	2:B:31:ALA:HB3	1.95	0.48
1:A:480:GLU:HA	1:A:481:PRO:HD3	1.75	0.48
1:A:666:THR:HB	1:A:689:ASN:HD22	1.78	0.48
1:A:569:ASP:O	1:A:571:GLU:N	2.35	0.48
1:A:655:LEU:O	1:A:657:ASN:N	2.47	0.47
1:A:430:LEU:HB3	1:A:434:GLY:HA2	1.95	0.47
1:A:695:PRO:HD2	1:A:698:LEU:HD12	1.95	0.47
1:A:541:MET:O	1:A:542:THR:C	2.53	0.47
1:A:440:ASN:O	1:A:443:ALA:HB3	2.15	0.47
1:A:675:THR:O	1:A:685:VAL:HA	2.14	0.47
2:B:85:TYR:O	2:B:86:ARG:C	2.52	0.47
1:A:685:VAL:HB	1:A:746:PHE:CE2	2.50	0.47
1:A:726:GLU:HG3	1:A:733:ILE:HG22	1.97	0.46
1:A:611:LEU:O	1:A:614:LEU:HB3	2.15	0.46
1:A:486:PHE:HB3	1:A:603:VAL:HA	1.97	0.46
1:A:541:MET:HG3	1:A:705:PRO:HG2	1.97	0.46
1:A:431:SER:O	1:A:433:ASP:N	2.49	0.46
1:A:648:ARG:O	1:A:652:LYS:HB2	2.16	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:724:ILE:O	1:A:728:GLU:HB2	2.16	0.46
1:A:569:ASP:C	1:A:571:GLU:H	2.16	0.46
1:A:468:ILE:HD11	1:A:487:TYR:CZ	2.50	0.46
1:A:618:VAL:HG12	1:A:622:PHE:HE2	1.80	0.46
1:A:440:ASN:OD1	1:A:441:LYS:N	2.49	0.45
1:A:569:ASP:CG	1:A:570:PRO:HD2	2.36	0.45
1:A:551:ASN:HB2	1:A:552:PRO:HD3	1.96	0.45
1:A:556:ILE:CD1	1:A:581:THR:HA	2.47	0.45
1:A:720:ILE:N	1:A:720:ILE:HD12	2.32	0.45
1:A:626:ILE:HG23	1:A:631:ILE:HB	1.98	0.45
1:A:454:VAL:HA	1:A:457:ARG:HE	1.81	0.45
1:A:481:PRO:HB3	1:A:500:PRO:HD2	1.98	0.45
1:A:520:THR:C	1:A:522:LEU:N	2.69	0.44
1:A:660:GLN:HA	1:A:663:ILE:HD12	1.98	0.44
1:A:679:MET:O	1:A:681:THR:N	2.50	0.44
1:A:538:LEU:O	1:A:542:THR:OG1	2.22	0.44
1:A:417:GLU:HG3	1:A:418:SER:N	2.33	0.44
2:B:83:SER:HB2	2:B:99:VAL:HA	2.00	0.44
1:A:700:GLU:HA	1:A:703:PHE:CD2	2.49	0.44
1:A:646:ALA:HB1	1:A:651:ILE:HD11	1.99	0.43
1:A:615:ILE:HG21	1:A:635:PHE:CZ	2.53	0.43
1:A:722:ARG:HA	1:A:733:ILE:HG21	1.99	0.43
1:A:626:ILE:HA	1:A:631:ILE:HD12	2.00	0.43
1:A:517:ASP:O	1:A:519:VAL:HG23	2.18	0.43
1:A:584:LEU:O	1:A:588:GLU:HG3	2.19	0.43
1:A:661:ASN:C	1:A:663:ILE:N	2.72	0.43
1:A:444:GLU:HG2	1:A:450:LYS:HA	2.00	0.43
1:A:650:ARG:NH2	2:B:85:TYR:HB2	2.34	0.42
1:A:648:ARG:C	1:A:650:ARG:N	2.72	0.42
1:A:666:THR:HB	1:A:689:ASN:HB2	2.00	0.42
1:A:629:MET:O	1:A:630:ASN:C	2.57	0.42
1:A:578:ASN:O	1:A:582:ASN:HB2	2.19	0.42
2:B:83:SER:OG	2:B:100:LYS:N	2.39	0.42
1:A:589:THR:HA	1:A:592:LYS:HE2	2.02	0.42
1:A:537:ILE:C	1:A:539:GLY:H	2.23	0.41
1:A:589:THR:HA	1:A:592:LYS:HG2	2.02	0.41
2:B:60:GLY:HA2	2:B:63:VAL:HB	2.02	0.41
1:A:691:GLY:HA3	1:A:692:PRO:HD3	1.81	0.41
1:A:585:SER:HB3	2:B:102:PHE:CE1	2.55	0.41
1:A:646:ALA:HB3	1:A:651:ILE:HD11	2.03	0.41
1:A:459:LEU:HG	1:A:465:PHE:CE1	2.56	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:633:PHE:CZ	1:A:674:ILE:CD1	2.93	0.41
1:A:651:ILE:HG12	1:A:729:HIS:NE2	2.35	0.41
1:A:589:THR:C	1:A:591:VAL:H	2.23	0.41
1:A:639:ASN:ND2	1:A:641:ASP:O	2.54	0.40
1:A:737:ASN:O	1:A:738:ARG:HB3	2.20	0.40
1:A:593:GLU:O	1:A:596:GLU:HB3	2.21	0.40
1:A:426:ALA:HA	1:A:440:ASN:ND2	2.35	0.40
1:A:596:GLU:O	1:A:598:SER:N	2.54	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	328/349 (94%)	238 (73%)	66 (20%)	24 (7%)	1	21
2	B	114/116 (98%)	98 (86%)	14 (12%)	2 (2%)	10	50
All	All	442/465 (95%)	336 (76%)	80 (18%)	26 (6%)	2	26

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	421	GLU
1	A	446	LEU
1	A	630	ASN
1	A	656	ILE
1	A	680	TYR
1	A	706	PHE
1	A	432	LYS
1	A	520	THR
1	A	623	GLU
1	A	649	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	86	ARG
1	A	521	GLU
1	A	570	PRO
1	A	464	ASP
1	A	479	LYS
1	A	486	PHE
2	B	25	GLY
1	A	416	SER
1	A	600	GLU
1	A	597	TYR
1	A	590	ILE
1	A	692	PRO
1	A	705	PRO
1	A	717	GLY
1	A	724	ILE
1	A	654	VAL

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	309/322 (96%)	266 (86%)	43 (14%)	4	27
2	B	103/100 (103%)	99 (96%)	4 (4%)	37	70
All	All	412/422 (98%)	365 (89%)	47 (11%)	7	34

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

Mol	Chain	Res	Type
1	A	427	ILE
1	A	431	SER
1	A	433	ASP
1	A	437	THR
1	A	438	GLU
1	A	455	LEU
1	A	461	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	474	SER
1	A	476	PHE
1	A	509	LEU
1	A	514	ILE
1	A	515	THR
1	A	517	ASP
1	A	520	THR
1	A	526	GLU
1	A	535	LEU
1	A	537	ILE
1	A	550	ARG
1	A	559	PHE
1	A	567	LEU
1	A	582	ASN
1	A	596	GLU
1	A	597	TYR
1	A	600	GLU
1	A	615	ILE
1	A	618	VAL
1	A	629	MET
1	A	632	ASP
1	A	634	CYS
1	A	644	VAL
1	A	651	ILE
1	A	653	GLN
1	A	664	GLU
1	A	674	ILE
1	A	686	SER
1	A	716	LEU
1	A	719	SER
1	A	723	LYS
1	A	732	LYS
1	A	734	TRP
1	A	736	GLU
1	A	737	ASN
1	A	742	VAL
2	B	48	LEU
2	B	82	TYR
2	B	89	MSE
2	B	105	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	495	ASN
1	A	547	HIS
1	A	669	ASN
1	A	729	HIS
1	A	737	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 1 ligands modelled in this entry, 1 is 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	334/349 (95%)	0.09	20 (5%) 23 16	161, 165, 165, 165	0
2	B	113/116 (97%)	0.61	20 (17%) 2 2	165, 165, 165, 165	0
All	All	447/465 (96%)	0.23	40 (8%) 10 8	161, 165, 165, 165	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	428	ILE	5.2
2	B	49	VAL	4.7
2	B	81	ALA	4.5
2	B	34	GLY	4.1
1	A	744	PHE	4.1
2	B	41	PHE	3.9
1	A	459	LEU	3.8
2	B	50	ILE	3.8
2	B	48	LEU	3.6
2	B	82	TYR	3.6
2	B	104	PHE	3.5
2	B	4	ILE	3.4
1	A	489	PHE	3.3
2	B	38	LEU	3.2
2	B	5	LEU	3.1
1	A	672	ILE	3.1
1	A	716	LEU	3.1
1	A	683	VAL	2.9
1	A	687	VAL	2.9
1	A	685	VAL	2.9
1	A	622	PHE	2.8
2	B	17	LEU	2.8
1	A	686	SER	2.7
1	A	468	ILE	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	427	ILE	2.7
2	B	107	LEU	2.7
2	B	6	VAL	2.5
2	B	51	LEU	2.4
1	A	516	ILE	2.4
1	A	514	ILE	2.3
1	A	462	LEU	2.3
2	B	58	ILE	2.2
1	A	556	ILE	2.2
2	B	84	HIS	2.2
1	A	563	MET	2.1
1	A	673	LYS	2.1
2	B	77	ILE	2.1
1	A	447	PHE	2.0
2	B	42	PHE	2.0
2	B	70	LYS	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(\AA^2)	Q<0.9
3	HG	A	800	1/1	0.96	0.04	-2.31	115,115,115,115	0

6.5 Other polymers [i](#)

There are no such residues in this entry.