Full wwPDB X-ray Structure Validation Report (i)

Feb 11, 2024 – 10:00 AM EST

PDB ID	:	3CIO
Title	:	The kinase domain of Escherichia coli tyrosine kinase ETK
Authors	:	Lee, D.C.; Zheng, J.; Jia, Z.
Deposited on	:	2008-03-11
Resolution	:	2.50 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:

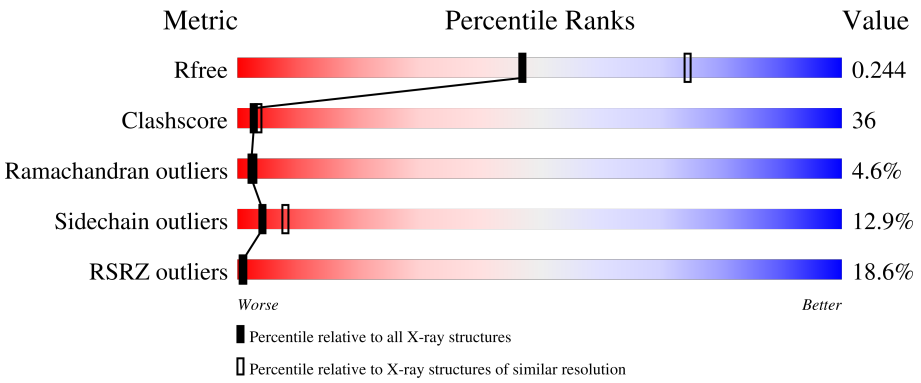
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	299	16%; 45%; 33%; 7%; 15%
1	D	299	16%; 40%; 36%; 9%; 15%

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	D	727	-	-	X	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4062 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 Tyrosine-protein kinase etk.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	255	1974	1240	349	377	8	0	0	0
1	D	255	1971	1238	350	375	8	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	428	MET	-	expression tag	UNP P38134
A	429	GLY	-	expression tag	UNP P38134
A	430	HIS	-	expression tag	UNP P38134
A	431	HIS	-	expression tag	UNP P38134
A	432	HIS	-	expression tag	UNP P38134
A	433	HIS	-	expression tag	UNP P38134
A	434	HIS	-	expression tag	UNP P38134
A	435	HIS	-	expression tag	UNP P38134
A	436	HIS	-	expression tag	UNP P38134
A	437	HIS	-	expression tag	UNP P38134
A	438	HIS	-	expression tag	UNP P38134
A	439	HIS	-	expression tag	UNP P38134
A	440	SER	-	expression tag	UNP P38134
A	441	SER	-	expression tag	UNP P38134
A	442	GLY	-	expression tag	UNP P38134
A	443	HIS	-	expression tag	UNP P38134
A	444	ILE	-	expression tag	UNP P38134
A	445	GLU	-	expression tag	UNP P38134
A	446	GLY	-	expression tag	UNP P38134
A	447	ARG	-	expression tag	UNP P38134
A	448	HIS	-	expression tag	UNP P38134
A	449	ILE	-	expression tag	UNP P38134
A	450	GLY	-	expression tag	UNP P38134
A	451	SER	-	expression tag	UNP P38134
D	428	MET	-	expression tag	UNP P38134

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Chain	Residue	Modelled	Actual	Comment	Reference
D	429	GLY	-	expression tag	UNP P38134
D	430	HIS	-	expression tag	UNP P38134
D	431	HIS	-	expression tag	UNP P38134
D	432	HIS	-	expression tag	UNP P38134
D	433	HIS	-	expression tag	UNP P38134
D	434	HIS	-	expression tag	UNP P38134
D	435	HIS	-	expression tag	UNP P38134
D	436	HIS	-	expression tag	UNP P38134
D	437	HIS	-	expression tag	UNP P38134
D	438	HIS	-	expression tag	UNP P38134
D	439	HIS	-	expression tag	UNP P38134
D	440	SER	-	expression tag	UNP P38134
D	441	SER	-	expression tag	UNP P38134
D	442	GLY	-	expression tag	UNP P38134
D	443	HIS	-	expression tag	UNP P38134
D	444	ILE	-	expression tag	UNP P38134
D	445	GLU	-	expression tag	UNP P38134
D	446	GLY	-	expression tag	UNP P38134
D	447	ARG	-	expression tag	UNP P38134
D	448	HIS	-	expression tag	UNP P38134
D	449	ILE	-	expression tag	UNP P38134
D	450	GLY	-	expression tag	UNP P38134
D	451	SER	-	expression tag	UNP P38134

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

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

- Molecule 3 is water.

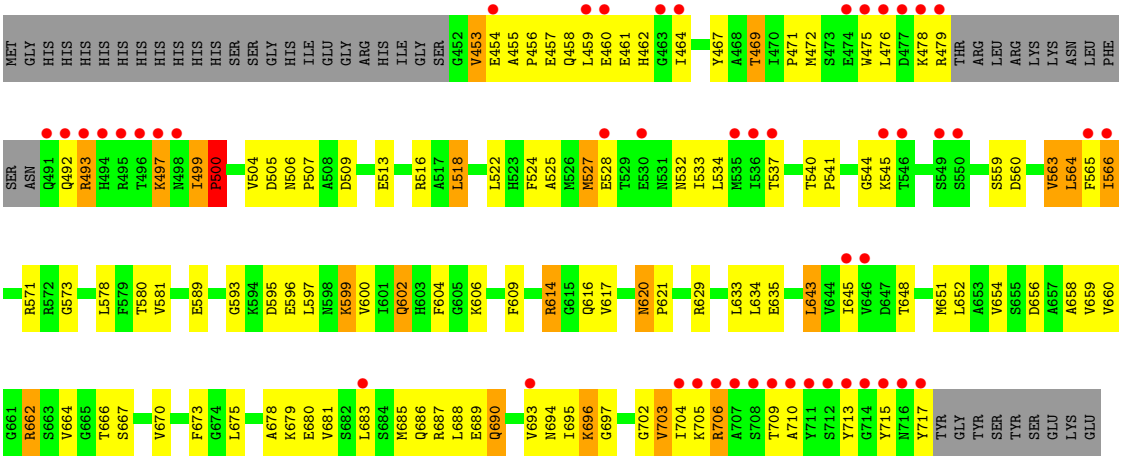
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	57	57	57	0	0
			Total	O		
3	D	57	57	57	0	0

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.

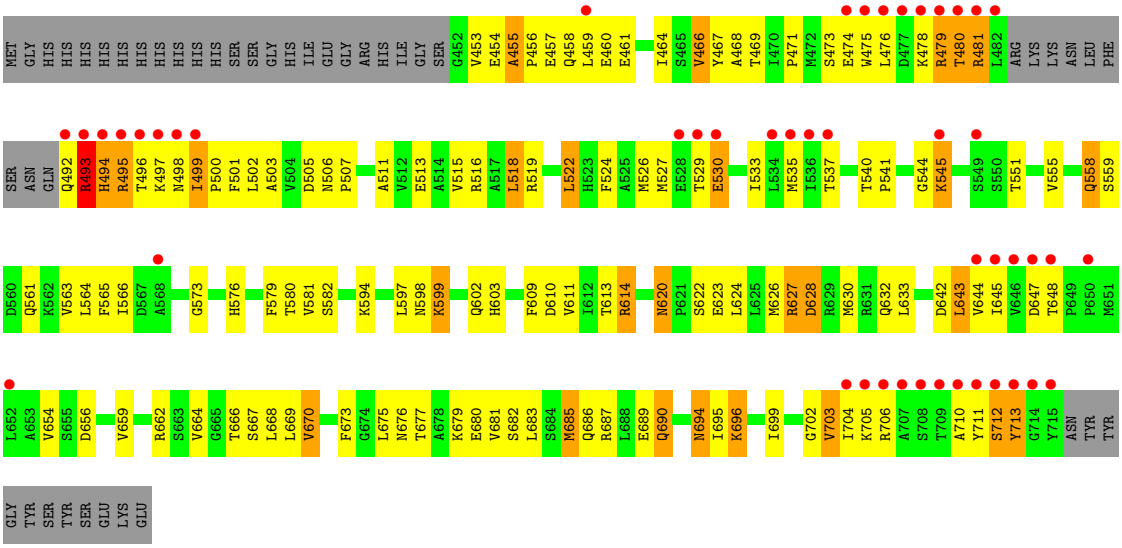
- Molecule 1: Tyrosine-protein kinase etk

Chain A:



- Molecule 1: Tyrosine-protein kinase etk

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	115.51Å 51.73Å 120.52Å 90.00° 114.57° 90.00°	Depositor
Resolution (Å)	29.93 – 2.50 29.93 – 2.50	Depositor EDS
% Data completeness (in resolution range)	69.2 (29.93-2.50) 94.1 (29.93-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	1.68 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.196 , 0.254 0.223 , 0.244	Depositor DCC
R_{free} test set	1095 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å^2)	51.2	Xtriage
Anisotropy	0.800	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 75.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4062	wwPDB-VP
Average B, all atoms (Å^2)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.61% 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)

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	0.39	0/2009	0.61	0/2722
1	D	0.39	0/2005	0.62	0/2716
All	All	0.39	0/4014	0.62	0/5438

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	1974	0	1965	112	0
1	D	1971	0	1973	171	0
2	A	2	0	0	0	0
2	D	1	0	0	3	0
3	A	57	0	0	3	0
3	D	57	0	0	12	0
All	All	4062	0	3938	284	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 36.

All (284) 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:D:476:LEU:HD11	1:D:502:LEU:HD13	1.29	1.14
1:D:705:LYS:HD3	1:D:706:ARG:H	1.00	1.12
1:D:499:ILE:HG12	1:D:500:PRO:HD2	1.32	1.11
1:A:453:VAL:HG13	1:A:454:GLU:H	1.27	0.99
1:D:620:ASN:HD22	1:D:620:ASN:H	1.07	0.99
1:D:495:ARG:H	1:D:495:ARG:HE	1.00	0.99
1:D:455:ALA:HB1	1:D:456:PRO:C	1.84	0.98
1:D:680:GLU:HA	1:D:683:LEU:HD12	1.48	0.94
1:D:545:LYS:HD3	1:D:545:LYS:H	1.32	0.93
1:D:705:LYS:HD3	1:D:706:ARG:N	1.84	0.93
1:A:620:ASN:HD22	1:A:620:ASN:H	1.16	0.93
1:A:472:MET:HG2	1:A:702:GLY:HA2	1.54	0.88
1:D:495:ARG:HE	1:D:495:ARG:N	1.72	0.86
1:D:456:PRO:HB3	1:D:466:VAL:HG11	1.59	0.84
1:D:495:ARG:H	1:D:495:ARG:NE	1.75	0.82
1:A:469:THR:HG23	1:A:703:VAL:HG21	1.61	0.81
1:A:499:ILE:HD13	1:A:499:ILE:H	1.45	0.81
1:D:476:LEU:HD21	1:D:502:LEU:HB2	1.63	0.80
1:D:662:ARG:HB2	1:D:662:ARG:NH2	1.97	0.80
1:D:545:LYS:HD3	1:D:545:LYS:N	1.95	0.79
1:D:705:LYS:CD	1:D:706:ARG:H	1.89	0.79
1:D:545:LYS:HG3	3:D:756:HOH:O	1.82	0.78
1:D:679:LYS:O	1:D:683:LEU:HG	1.84	0.78
1:D:694:ASN:HB3	1:D:696:LYS:HE2	1.64	0.77
1:A:492:GLN:O	1:A:493:ARG:HG2	1.85	0.77
1:A:620:ASN:HD22	1:A:620:ASN:N	1.82	0.76
1:D:623:GLU:HB3	3:D:766:HOH:O	1.85	0.76
1:A:457:GLU:HA	1:A:460:GLU:HG3	1.67	0.76
1:A:479:ARG:HB2	1:A:479:ARG:HH21	1.51	0.76
1:A:545:LYS:HE3	1:A:648:THR:O	1.87	0.74
1:A:679:LYS:HG2	1:A:683:LEU:HD11	1.67	0.74
1:D:573:GLY:O	1:D:614:ARG:HD3	1.88	0.74
1:D:620:ASN:H	1:D:620:ASN:ND2	1.85	0.74
1:D:499:ILE:CG1	1:D:500:PRO:HD2	2.15	0.73
1:D:679:LYS:HG2	1:D:683:LEU:HD11	1.71	0.72
1:A:686:GLN:HE21	1:A:690:GLN:HE22	1.38	0.72
1:D:533:ILE:HG22	1:D:664:VAL:HG23	1.71	0.71
1:D:656:ASP:O	1:D:659:VAL:HG12	1.91	0.71
1:D:662:ARG:CB	1:D:662:ARG:HH21	2.04	0.71
1:D:455:ALA:HB2	1:D:458:GLN:HB3	1.72	0.70
1:D:518:LEU:HD22	1:D:522:LEU:HD23	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:478:LYS:O	1:D:479:ARG:HB2	1.90	0.69
1:D:505:ASP:O	1:D:507:PRO:HD3	1.93	0.69
1:D:597:LEU:HD11	1:D:633:LEU:HD13	1.75	0.69
1:D:669:LEU:HD23	1:D:681:VAL:HG23	1.74	0.69
1:D:673:PHE:CD1	1:D:703:VAL:HA	2.28	0.69
1:A:532:ASN:HB2	3:A:753:HOH:O	1.92	0.68
1:D:624:LEU:O	1:D:627:ARG:HB2	1.94	0.68
1:A:479:ARG:HB2	1:A:479:ARG:NH2	2.08	0.68
1:A:656:ASP:O	1:A:660:VAL:HG22	1.93	0.67
1:D:524:PHE:O	1:D:527:MET:HB2	1.94	0.67
1:D:581:VAL:HG22	1:D:602:GLN:NE2	2.09	0.67
1:A:571:ARG:HD2	3:A:748:HOH:O	1.95	0.67
1:D:620:ASN:HD22	1:D:620:ASN:N	1.77	0.66
1:A:573:GLY:O	1:A:614:ARG:HD3	1.95	0.66
1:D:662:ARG:HB2	1:D:662:ARG:HH21	1.55	0.66
1:D:474:GLU:HG3	1:D:706:ARG:NH2	2.09	0.66
1:A:652:LEU:HD11	1:A:680:GLU:HB2	1.78	0.66
1:D:455:ALA:HB3	1:D:459:LEU:HD23	1.78	0.66
1:D:493:ARG:HA	1:D:493:ARG:HH11	1.59	0.65
1:A:705:LYS:HG2	1:A:706:ARG:H	1.61	0.65
1:D:473:SER:OG	1:D:476:LEU:HD13	1.97	0.64
1:D:696:LYS:HG2	3:D:757:HOH:O	1.98	0.64
1:A:453:VAL:HG13	1:A:454:GLU:N	2.06	0.64
1:A:678:ALA:O	1:A:681:VAL:HG22	1.98	0.64
1:D:558:GLN:HE21	1:D:558:GLN:N	1.95	0.64
1:A:620:ASN:N	1:A:620:ASN:ND2	2.47	0.63
1:D:515:VAL:HG21	1:D:551:THR:HG21	1.81	0.63
1:A:524:PHE:O	1:A:527:MET:HB2	1.99	0.62
1:A:499:ILE:H	1:A:499:ILE:CD1	2.12	0.62
1:A:497:LYS:HB2	1:A:497:LYS:NZ	2.15	0.62
1:A:662:ARG:HH21	1:A:662:ARG:HB2	1.63	0.62
1:D:581:VAL:HG22	1:D:602:GLN:HE22	1.65	0.62
1:D:681:VAL:HG13	1:D:682:SER:N	2.15	0.61
1:D:467:TYR:OH	1:D:696:LYS:HB3	2.01	0.60
1:D:522:LEU:HD12	1:D:526:MET:SD	2.41	0.60
1:A:620:ASN:H	1:A:620:ASN:ND2	1.93	0.60
1:A:559:SER:O	1:A:560:ASP:HB2	2.02	0.60
1:D:541:PRO:HD3	3:D:749:HOH:O	2.01	0.60
1:D:581:VAL:CG2	1:D:602:GLN:NE2	2.64	0.60
1:D:623:GLU:OE1	1:D:623:GLU:N	2.33	0.60
1:D:681:VAL:HG13	1:D:682:SER:H	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:603:HIS:HD2	1:D:610:ASP:OD2	1.84	0.59
1:D:455:ALA:CB	1:D:456:PRO:CA	2.80	0.59
1:A:544:GLY:HA2	3:A:730:HOH:O	2.01	0.59
1:D:527:MET:O	1:D:529:THR:HG23	2.02	0.59
1:A:643:LEU:HD13	1:A:645:ILE:HG13	1.85	0.59
1:D:686:GLN:HG2	1:D:690:GLN:HE22	1.67	0.59
1:A:499:ILE:HD13	1:A:499:ILE:N	2.16	0.59
1:D:576:HIS:HB3	1:D:613:THR:CG2	2.32	0.58
1:D:540:THR:HA	3:D:749:HOH:O	2.04	0.58
1:D:455:ALA:CB	1:D:456:PRO:C	2.68	0.58
1:D:597:LEU:HD11	1:D:633:LEU:CD1	2.32	0.58
1:A:504:VAL:HG21	1:A:606:LYS:HG2	1.86	0.58
1:D:474:GLU:HG3	1:D:706:ARG:HH21	1.66	0.58
1:A:616:GLN:HA	1:A:616:GLN:NE2	2.19	0.58
1:D:475:TRP:CH2	1:D:480:THR:HG21	2.39	0.57
1:D:473:SER:HA	3:D:765:HOH:O	2.05	0.57
1:A:476:LEU:HD13	1:A:476:LEU:O	2.04	0.57
1:D:545:LYS:NZ	2:D:727:CL:CL	2.72	0.57
1:D:526:MET:HE3	3:D:728:HOH:O	2.03	0.56
1:A:454:GLU:HA	1:A:673:PHE:CD2	2.40	0.56
1:D:457:GLU:CD	1:D:457:GLU:H	2.09	0.56
1:A:651:MET:HG3	1:A:688:LEU:HG	1.88	0.56
1:D:659:VAL:HG23	1:D:662:ARG:NH1	2.21	0.56
1:D:454:GLU:HB2	1:D:673:PHE:HE2	1.71	0.56
1:D:696:LYS:N	1:D:696:LYS:HD2	2.20	0.56
1:D:533:ILE:CG2	1:D:664:VAL:HG23	2.35	0.55
1:D:464:ILE:HD11	1:D:682:SER:HB3	1.88	0.55
1:D:522:LEU:HD12	1:D:526:MET:HG3	1.89	0.55
1:A:472:MET:HE2	1:A:472:MET:HA	1.88	0.55
1:A:679:LYS:O	1:A:683:LEU:HG	2.07	0.55
1:D:453:VAL:HG12	1:D:676:ASN:O	2.07	0.55
1:D:704:ILE:O	1:D:704:ILE:HG23	2.06	0.55
1:D:668:LEU:HD13	1:D:699:ILE:HD11	1.89	0.55
1:D:537:THR:HG23	1:D:669:LEU:HD12	1.89	0.54
1:A:581:VAL:HG22	1:A:602:GLN:NE2	2.23	0.54
1:D:659:VAL:HG23	1:D:662:ARG:HH11	1.72	0.54
1:D:576:HIS:HB3	1:D:613:THR:HG21	1.90	0.54
1:A:706:ARG:NH1	1:A:706:ARG:HB3	2.22	0.54
1:D:515:VAL:HG21	1:D:551:THR:CG2	2.37	0.54
1:D:711:TYR:O	1:D:712:SER:HB2	2.08	0.54
1:D:498:ASN:HA	1:D:580:THR:OG1	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:658:ALA:O	1:A:662:ARG:HG3	2.08	0.53
1:D:545:LYS:HD2	3:D:754:HOH:O	2.07	0.53
1:A:563:VAL:HG22	1:A:609:PHE:HB2	1.91	0.53
1:A:666:THR:HG23	1:A:696:LYS:HG2	1.90	0.53
1:A:499:ILE:HB	1:A:500:PRO:HB3	1.89	0.53
1:D:455:ALA:HB1	1:D:456:PRO:CA	2.38	0.53
1:D:675:LEU:HD13	1:D:675:LEU:C	2.28	0.53
1:D:505:ASP:C	1:D:507:PRO:HD3	2.29	0.53
1:A:687:ARG:HH21	1:A:687:ARG:HG3	1.71	0.53
1:A:673:PHE:CD2	1:A:703:VAL:HA	2.44	0.53
1:D:513:GLU:OE2	1:D:516:ARG:NH2	2.42	0.53
1:A:702:GLY:O	1:A:703:VAL:C	2.48	0.52
1:A:518:LEU:HD22	1:A:522:LEU:HD23	1.90	0.52
1:A:662:ARG:HH21	1:A:662:ARG:CB	2.21	0.52
1:D:563:VAL:HG22	1:D:609:PHE:HB2	1.92	0.52
1:A:497:LYS:HB2	1:A:497:LYS:HZ3	1.74	0.52
1:A:518:LEU:HD22	1:A:522:LEU:CD2	2.40	0.52
1:A:513:GLU:HA	1:A:516:ARG:NH2	2.25	0.52
1:D:643:LEU:HD13	1:D:645:ILE:HG13	1.90	0.52
1:A:533:ILE:O	1:A:664:VAL:HG13	2.10	0.51
1:A:564:LEU:HD22	1:A:565:PHE:N	2.25	0.51
1:A:537:THR:HB	1:A:648:THR:OG1	2.11	0.51
1:A:705:LYS:HG2	1:A:706:ARG:N	2.26	0.51
1:A:505:ASP:O	1:A:507:PRO:HD3	2.11	0.51
1:D:599:LYS:HB2	1:D:599:LYS:HZ3	1.76	0.51
1:A:593:GLY:HA2	1:A:629:ARG:NH1	2.26	0.50
1:A:597:LEU:O	1:A:600:VAL:HG22	2.12	0.50
1:A:599:LYS:HB2	1:A:599:LYS:NZ	2.26	0.50
1:D:666:THR:HG23	1:D:696:LYS:HD3	1.93	0.49
1:D:498:ASN:ND2	1:D:499:ILE:H	2.10	0.49
1:A:578:LEU:HD23	1:A:578:LEU:C	2.33	0.49
1:D:544:GLY:HA2	3:D:730:HOH:O	2.11	0.49
1:A:578:LEU:HD23	1:A:578:LEU:O	2.13	0.49
1:D:696:LYS:HD2	1:D:696:LYS:H	1.78	0.49
1:D:455:ALA:HB1	1:D:457:GLU:N	2.23	0.49
1:D:511:ALA:O	1:D:515:VAL:HG23	2.13	0.49
1:D:712:SER:O	1:D:713:TYR:HB2	2.13	0.49
1:A:479:ARG:HH21	1:A:479:ARG:CB	2.22	0.48
1:A:455:ALA:O	1:A:458:GLN:HB3	2.13	0.48
1:D:476:LEU:HD11	1:D:502:LEU:CD1	2.21	0.48
1:A:685:MET:O	1:A:689:GLU:HG2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:GLU:OE2	1:A:457:GLU:N	2.44	0.48
1:D:537:THR:HB	1:D:648:THR:OG1	2.13	0.48
1:D:705:LYS:O	1:D:706:ARG:HG2	2.13	0.48
1:A:471:PRO:O	1:A:472:MET:HE2	2.13	0.48
1:A:472:MET:HA	1:A:472:MET:CE	2.41	0.48
1:A:581:VAL:CG2	1:A:602:GLN:NE2	2.77	0.48
1:A:662:ARG:HB2	1:A:662:ARG:NH2	2.29	0.48
1:D:494:HIS:HA	1:D:495:ARG:HH11	1.79	0.48
1:D:623:GLU:H	1:D:623:GLU:CD	2.17	0.48
1:D:687:ARG:HG2	1:D:687:ARG:HH21	1.78	0.48
1:D:455:ALA:HB1	1:D:458:GLN:N	2.28	0.47
1:D:455:ALA:CB	1:D:458:GLN:H	2.27	0.47
1:A:666:THR:HG22	1:A:667:SER:N	2.29	0.47
1:A:467:TYR:CE2	1:A:697:GLY:HA3	2.49	0.47
1:A:597:LEU:HD11	1:A:633:LEU:HD13	1.96	0.47
1:D:694:ASN:HB3	1:D:696:LYS:CE	2.39	0.47
1:D:667:SER:HB2	1:D:695:ILE:HA	1.96	0.46
1:A:666:THR:HG22	1:A:667:SER:H	1.80	0.46
1:D:643:LEU:HD22	1:D:644:VAL:N	2.31	0.46
1:A:457:GLU:HA	1:A:460:GLU:CG	2.43	0.46
1:A:475:TRP:HA	1:A:478:LYS:HD3	1.96	0.46
1:D:454:GLU:HB2	1:D:673:PHE:CE2	2.50	0.46
1:D:667:SER:HB2	1:D:694:ASN:O	2.16	0.46
1:D:476:LEU:HD12	1:D:476:LEU:N	2.31	0.46
1:D:622:SER:HB2	1:D:623:GLU:OE1	2.16	0.46
1:D:480:THR:O	1:D:481:ARG:HB3	2.15	0.46
1:A:540:THR:HB	1:A:541:PRO:CD	2.46	0.45
1:A:664:VAL:HG12	1:A:666:THR:O	2.16	0.45
1:D:516:ARG:O	1:D:519:ARG:HB3	2.16	0.45
1:D:474:GLU:O	1:D:478:LYS:HG3	2.16	0.45
1:D:479:ARG:O	1:D:480:THR:HG22	2.16	0.45
1:D:535:MET:N	1:D:664:VAL:HG21	2.32	0.45
1:A:564:LEU:HD13	1:A:566:ILE:HG22	1.99	0.45
1:A:688:LEU:HD12	1:A:695:ILE:HD11	1.98	0.45
1:A:563:VAL:HG22	1:A:609:PHE:CB	2.46	0.45
1:A:688:LEU:CD1	1:A:695:ILE:HD11	2.46	0.45
1:A:493:ARG:HG2	1:A:493:ARG:HH21	1.81	0.45
1:A:614:ARG:HH11	1:A:617:VAL:HG23	1.81	0.45
1:A:706:ARG:CB	1:A:706:ARG:HH11	2.30	0.45
1:D:581:VAL:HG22	1:D:582:SER:H	1.81	0.45
1:A:656:ASP:O	1:A:659:VAL:HG22	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:597:LEU:CD1	1:D:633:LEU:HD13	2.45	0.44
1:D:469:THR:CG2	1:D:703:VAL:HG21	2.47	0.44
1:D:675:LEU:HD13	1:D:675:LEU:O	2.17	0.44
1:D:499:ILE:HG23	1:D:500:PRO:N	2.32	0.44
1:D:468:ALA:HA	3:D:776:HOH:O	2.16	0.44
1:A:464:ILE:CD1	1:A:681:VAL:HG23	2.48	0.44
1:A:571:ARG:CZ	1:A:621:PRO:HB2	2.48	0.44
1:D:594:LYS:HA	3:D:732:HOH:O	2.18	0.44
1:D:453:VAL:HG13	1:D:453:VAL:O	2.18	0.43
1:D:499:ILE:HG23	1:D:500:PRO:CD	2.48	0.43
1:A:455:ALA:HA	1:A:456:PRO:HD3	1.94	0.43
1:D:599:LYS:HB2	1:D:599:LYS:NZ	2.31	0.43
1:D:626:MET:HG2	1:D:659:VAL:HG21	2.00	0.43
1:D:630:MET:HE3	1:D:630:MET:HA	2.00	0.43
1:A:455:ALA:HB3	1:A:458:GLN:HE21	1.83	0.43
1:D:540:THR:HB	1:D:541:PRO:HD2	2.00	0.43
1:A:688:LEU:O	1:A:693:VAL:HB	2.18	0.43
1:D:493:ARG:NH1	1:D:495:ARG:HD3	2.33	0.43
1:A:513:GLU:HA	1:A:516:ARG:HH21	1.83	0.43
1:A:589:GLU:HB3	1:A:595:ASP:OD2	2.18	0.43
1:D:498:ASN:HD22	1:D:499:ILE:H	1.66	0.43
1:A:459:LEU:N	1:A:459:LEU:HD22	2.34	0.43
1:D:662:ARG:NH2	1:D:662:ARG:CB	2.68	0.43
1:D:687:ARG:HG2	1:D:687:ARG:NH2	2.34	0.43
1:D:522:LEU:HD13	1:D:522:LEU:HA	1.81	0.43
1:D:492:GLN:N	1:D:492:GLN:OE1	2.51	0.42
1:D:627:ARG:O	1:D:628:ASP:HB2	2.17	0.42
1:D:677:THR:C	1:D:679:LYS:N	2.72	0.42
1:A:497:LYS:CD	1:A:580:THR:HB	2.49	0.42
1:D:522:LEU:HD12	1:D:526:MET:CG	2.49	0.42
1:D:702:GLY:O	1:D:703:VAL:C	2.57	0.42
1:D:469:THR:HG23	1:D:703:VAL:HG21	2.01	0.42
1:A:497:LYS:HD3	1:A:580:THR:HB	2.01	0.42
1:D:471:PRO:HA	1:D:703:VAL:O	2.19	0.42
1:D:475:TRP:CZ2	1:D:480:THR:HG21	2.54	0.42
1:A:509:ASP:O	1:A:513:GLU:HG2	2.20	0.42
1:A:518:LEU:O	1:A:522:LEU:HD23	2.20	0.42
1:A:524:PHE:O	1:A:525:ALA:C	2.57	0.42
1:D:545:LYS:CE	2:D:727:CL:CL	3.05	0.42
1:A:687:ARG:HD3	1:A:687:ARG:HA	1.89	0.42
1:A:680:GLU:HA	1:A:683:LEU:HD12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:704:ILE:O	1:A:704:ILE:CG2	2.67	0.42
1:D:677:THR:C	1:D:679:LYS:H	2.23	0.42
1:A:504:VAL:HG11	1:A:606:LYS:O	2.20	0.42
1:D:581:VAL:HG22	1:D:582:SER:N	2.34	0.42
1:A:534:LEU:CD1	1:A:666:THR:HB	2.50	0.41
1:D:460:GLU:OE1	1:D:466:VAL:HG12	2.20	0.41
1:D:673:PHE:CE1	1:D:704:ILE:HG22	2.55	0.41
1:A:533:ILE:HD13	1:A:634:LEU:HB3	2.02	0.41
1:A:685:MET:HG2	1:A:695:ILE:HG21	2.01	0.41
1:D:558:GLN:HE21	1:D:558:GLN:H	1.64	0.41
1:D:620:ASN:ND2	1:D:620:ASN:N	2.45	0.41
1:A:497:LYS:HE2	1:A:604:PHE:CE2	2.56	0.41
1:D:455:ALA:HB1	1:D:458:GLN:H	1.85	0.41
1:D:475:TRP:HE3	1:D:476:LEU:HD12	1.84	0.41
1:D:493:ARG:CZ	1:D:495:ARG:HD3	2.50	0.41
1:D:537:THR:O	1:D:670:VAL:HG12	2.21	0.41
1:D:656:ASP:HA	1:D:659:VAL:HG12	2.02	0.41
1:A:534:LEU:HD13	1:A:666:THR:HB	2.02	0.41
1:D:501:PHE:CE2	1:D:503:ALA:HB3	2.55	0.41
1:D:565:PHE:HB3	1:D:611:VAL:HG12	2.02	0.41
1:D:579:PHE:O	1:D:580:THR:C	2.59	0.41
1:D:614:ARG:HE	1:D:614:ARG:HB2	1.62	0.41
2:D:727:CL:CL	3:D:743:HOH:O	2.60	0.41
1:D:497:LYS:O	1:D:497:LYS:HG2	2.19	0.41
1:D:454:GLU:H	1:D:458:GLN:NE2	2.19	0.41
1:D:464:ILE:HG23	1:D:685:MET:HE1	2.02	0.41
1:D:551:THR:O	1:D:555:VAL:HG23	2.21	0.41
1:A:532:ASN:HD22	1:A:532:ASN:HA	1.67	0.41
1:D:545:LYS:HA	1:D:670:VAL:HG11	2.03	0.41
1:D:711:TYR:HB3	1:D:712:SER:H	1.65	0.41
1:A:596:GLU:HA	1:A:596:GLU:OE1	2.21	0.41
1:D:594:LYS:HE2	1:D:594:LYS:HB3	1.78	0.41
1:D:476:LEU:O	1:D:479:ARG:N	2.54	0.40
1:D:676:ASN:HA	1:D:680:GLU:OE2	2.21	0.40
1:A:706:ARG:HB3	1:A:706:ARG:HH11	1.87	0.40
1:A:599:LYS:HB2	1:A:599:LYS:HZ3	1.86	0.40
1:D:681:VAL:CG1	1:D:682:SER:N	2.84	0.40
1:D:561:GLN:HG3	1:D:642:ASP:HB2	2.04	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	251/299 (84%)	221 (88%)	21 (8%)	9 (4%)	3 4
1	D	251/299 (84%)	212 (84%)	25 (10%)	14 (6%)	2 1
All	All	502/598 (84%)	433 (86%)	46 (9%)	23 (5%)	2 2

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	500	PRO
1	A	703	VAL
1	D	455	ALA
1	D	480	THR
1	D	481	ARG
1	D	627	ARG
1	D	628	ASP
1	A	706	ARG
1	A	710	ALA
1	D	493	ARG
1	D	694	ASN
1	D	703	VAL
1	A	493	ARG
1	A	696	LYS
1	D	530	GLU
1	D	710	ALA
1	D	712	SER
1	A	453	VAL
1	A	709	THR
1	D	496	THR
1	D	494	HIS
1	D	713	TYR
1	A	528	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	213/252 (84%)	186 (87%)	27 (13%)	4 8
1	D	213/252 (84%)	185 (87%)	28 (13%)	4 7
All	All	426/504 (84%)	371 (87%)	55 (13%)	4 8

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	461	GLU
1	A	462	HIS
1	A	469	THR
1	A	497	LYS
1	A	499	ILE
1	A	500	PRO
1	A	506	ASN
1	A	518	LEU
1	A	527	MET
1	A	563	VAL
1	A	564	LEU
1	A	566	ILE
1	A	599	LYS
1	A	602	GLN
1	A	614	ARG
1	A	620	ASN
1	A	635	GLU
1	A	643	LEU
1	A	654	VAL
1	A	662	ARG
1	A	670	VAL
1	A	675	LEU
1	A	690	GLN
1	A	694	ASN
1	A	713	TYR
1	A	715	TYR
1	A	717	TYR

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Mol	Chain	Res	Type
1	D	461	GLU
1	D	466	VAL
1	D	479	ARG
1	D	493	ARG
1	D	495	ARG
1	D	499	ILE
1	D	506	ASN
1	D	518	LEU
1	D	522	LEU
1	D	530	GLU
1	D	545	LYS
1	D	558	GLN
1	D	559	SER
1	D	564	LEU
1	D	566	ILE
1	D	598	ASN
1	D	599	LYS
1	D	614	ARG
1	D	620	ASN
1	D	632	GLN
1	D	643	LEU
1	D	647	ASP
1	D	654	VAL
1	D	670	VAL
1	D	685	MET
1	D	689	GLU
1	D	690	GLN
1	D	696	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	458	GLN
1	A	506	ASN
1	A	532	ASN
1	A	558	GLN
1	A	577	ASN
1	A	602	GLN
1	A	603	HIS
1	A	616	GLN
1	A	620	ASN
1	A	686	GLN

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Mol	Chain	Res	Type
1	D	458	GLN
1	D	506	ASN
1	D	532	ASN
1	D	558	GLN
1	D	577	ASN
1	D	602	GLN
1	D	603	HIS
1	D	620	ASN
1	D	640	HIS
1	D	694	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 3 ligands modelled in this entry, 3 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 (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	255/299 (85%)	1.19	48 (18%)	1	1	30, 64, 100, 100	0
1	D	255/299 (85%)	1.15	47 (18%)	1	1	32, 62, 100, 100	0
All	All	510/598 (85%)	1.17	95 (18%)	1	1	30, 63, 100, 100	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	710	ALA	12.0
1	D	708	SER	11.6
1	D	482	LEU	10.8
1	D	711	TYR	10.7
1	D	498	ASN	10.4
1	A	496	THR	10.0
1	D	709	THR	10.0
1	A	497	LYS	9.9
1	A	711	TYR	9.6
1	A	713	TYR	9.0
1	A	708	SER	8.4
1	D	707	ALA	8.1
1	A	493	ARG	7.6
1	A	709	THR	7.5
1	D	713	TYR	7.4
1	A	707	ALA	7.3
1	D	481	ARG	7.0
1	A	495	ARG	7.0
1	A	494	HIS	7.0
1	A	706	ARG	6.9
1	A	528	GLU	6.8
1	D	706	ARG	6.6
1	D	710	ALA	6.6
1	A	476	LEU	6.5

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Mol	Chain	Res	Type	RSRZ
1	D	494	HIS	6.4
1	A	714	GLY	5.8
1	A	715	TYR	5.6
1	D	495	ARG	5.6
1	D	704	ILE	5.6
1	A	705	LYS	5.6
1	A	491	GLN	5.6
1	D	712	SER	5.6
1	A	712	SER	5.4
1	D	480	THR	5.3
1	D	499	ILE	5.2
1	A	474	GLU	5.2
1	D	493	ARG	5.1
1	D	497	LYS	5.0
1	A	498	ASN	5.0
1	D	496	THR	5.0
1	D	476	LEU	4.9
1	A	492	GLN	4.9
1	A	463	GLY	4.9
1	D	479	ARG	4.9
1	D	705	LYS	4.7
1	A	464	ILE	4.6
1	D	714	GLY	4.4
1	A	704	ILE	4.4
1	D	492	GLN	4.3
1	D	459	LEU	4.3
1	A	536	ILE	4.2
1	A	475	TRP	4.1
1	A	716	ASN	4.1
1	D	645	ILE	3.8
1	D	536	ILE	3.7
1	D	478	LYS	3.6
1	A	645	ILE	3.5
1	D	528	GLU	3.4
1	A	459	LEU	3.2
1	A	478	LYS	3.2
1	D	474	GLU	3.0
1	A	477	ASP	2.8
1	D	534	LEU	2.8
1	A	549	SER	2.8
1	D	715	TYR	2.8
1	D	475	TRP	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	535	MET	2.8
1	A	545	LYS	2.6
1	A	546	THR	2.6
1	A	535	MET	2.6
1	D	477	ASP	2.6
1	A	460	GLU	2.5
1	A	454	GLU	2.5
1	D	530	GLU	2.5
1	D	652	LEU	2.4
1	A	537	THR	2.4
1	A	646	VAL	2.4
1	D	537	THR	2.3
1	D	646	VAL	2.3
1	D	644	VAL	2.3
1	D	568	ALA	2.3
1	D	549	SER	2.2
1	A	479	ARG	2.2
1	D	648	THR	2.2
1	D	650	PRO	2.2
1	A	566	ILE	2.1
1	A	717	TYR	2.1
1	A	530	GLU	2.1
1	D	647	ASP	2.1
1	D	529	THR	2.1
1	D	545	LYS	2.1
1	A	693	VAL	2.1
1	A	565	PHE	2.1
1	A	683	LEU	2.1
1	A	550	SER	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)

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	D	727	1/1	0.57	0.23	100,100,100,100	0
2	CL	A	728	1/1	0.92	0.42	75,75,75,75	0
2	CL	A	727	1/1	0.92	0.33	60,60,60,60	0

6.5 Other polymers (i)

There are no such residues in this entry.