Full wwPDB X-ray Structure Validation Report (i)

Nov 15, 2023 – 10:08 PM JST

PDB ID	:	6K0K
Title	:	Crystal structure of Escherichia coli pyruvate kinase II
Authors	:	Zhao, C.H.; Zhang, Y.P.; Li, Y.
Deposited on	:	2019-05-07
Resolution	:	2.68 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 (i)) were used in the production of this report:

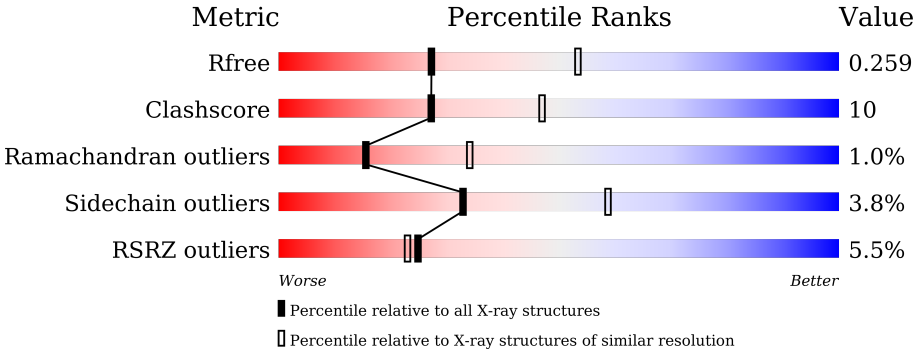
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.68 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	3863 (2.70-2.66)
Clashscore	141614	4210 (2.70-2.66)
Ramachandran outliers	138981	4141 (2.70-2.66)
Sidechain outliers	138945	4141 (2.70-2.66)
RSRZ outliers	127900	3780 (2.70-2.66)

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	489	7%, 74%, 22%, • •
1	B	489	4%, 78%, 18%, • •
1	C	489	4%, 61%, 19%, •, 19%
1	D	489	5%, 56%, 23%, •, 20%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CIT	A	502	-	X	-	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 13318 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 Pyruvate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	479	3579	2220	643	693	23	0	0	0
1	B	478	3573	2217	642	691	23	0	0	0
1	C	397	2955	1820	541	572	22	0	0	0
1	D	392	2924	1801	536	565	22	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	481	ALA	-	expression tag	UNP C3T5U7
A	482	LEU	-	expression tag	UNP C3T5U7
A	483	GLY	-	expression tag	UNP C3T5U7
A	484	HIS	-	expression tag	UNP C3T5U7
A	485	HIS	-	expression tag	UNP C3T5U7
A	486	HIS	-	expression tag	UNP C3T5U7
A	487	HIS	-	expression tag	UNP C3T5U7
A	488	HIS	-	expression tag	UNP C3T5U7
A	489	HIS	-	expression tag	UNP C3T5U7
B	481	ALA	-	expression tag	UNP C3T5U7
B	482	LEU	-	expression tag	UNP C3T5U7
B	483	GLY	-	expression tag	UNP C3T5U7
B	484	HIS	-	expression tag	UNP C3T5U7
B	485	HIS	-	expression tag	UNP C3T5U7
B	486	HIS	-	expression tag	UNP C3T5U7
B	487	HIS	-	expression tag	UNP C3T5U7
B	488	HIS	-	expression tag	UNP C3T5U7
B	489	HIS	-	expression tag	UNP C3T5U7
C	481	ALA	-	expression tag	UNP C3T5U7
C	482	LEU	-	expression tag	UNP C3T5U7
C	483	GLY	-	expression tag	UNP C3T5U7

Continued on next page...

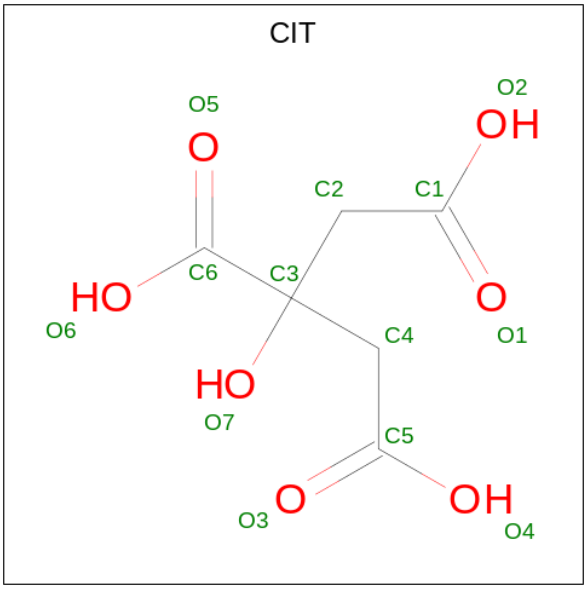
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	484	HIS	-	expression tag	UNP C3T5U7
C	485	HIS	-	expression tag	UNP C3T5U7
C	486	HIS	-	expression tag	UNP C3T5U7
C	487	HIS	-	expression tag	UNP C3T5U7
C	488	HIS	-	expression tag	UNP C3T5U7
C	489	HIS	-	expression tag	UNP C3T5U7
D	481	ALA	-	expression tag	UNP C3T5U7
D	482	LEU	-	expression tag	UNP C3T5U7
D	483	GLY	-	expression tag	UNP C3T5U7
D	484	HIS	-	expression tag	UNP C3T5U7
D	485	HIS	-	expression tag	UNP C3T5U7
D	486	HIS	-	expression tag	UNP C3T5U7
D	487	HIS	-	expression tag	UNP C3T5U7
D	488	HIS	-	expression tag	UNP C3T5U7
D	489	HIS	-	expression tag	UNP C3T5U7

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	13	6	7	0	0
			Total	C	O		
3	C	1	13	6	7	0	0
			Total	C	O		
3	D	1	13	6	7	0	0

- Molecule 4 is water.

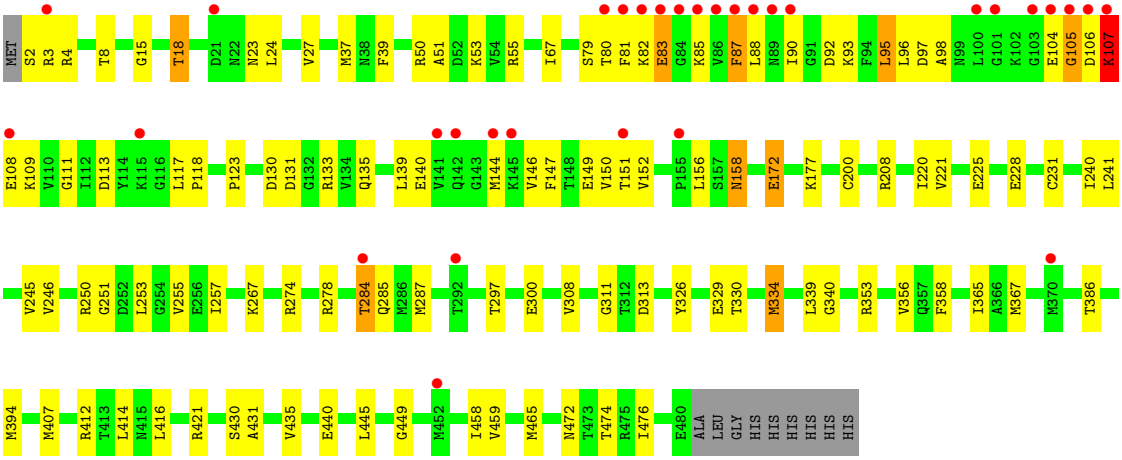
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	77	77	77	0	0
			Total	O		
4	B	105	105	105	0	0
			Total	O		
4	C	47	47	47	0	0
			Total	O		
4	D	15	15	15	0	0

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.

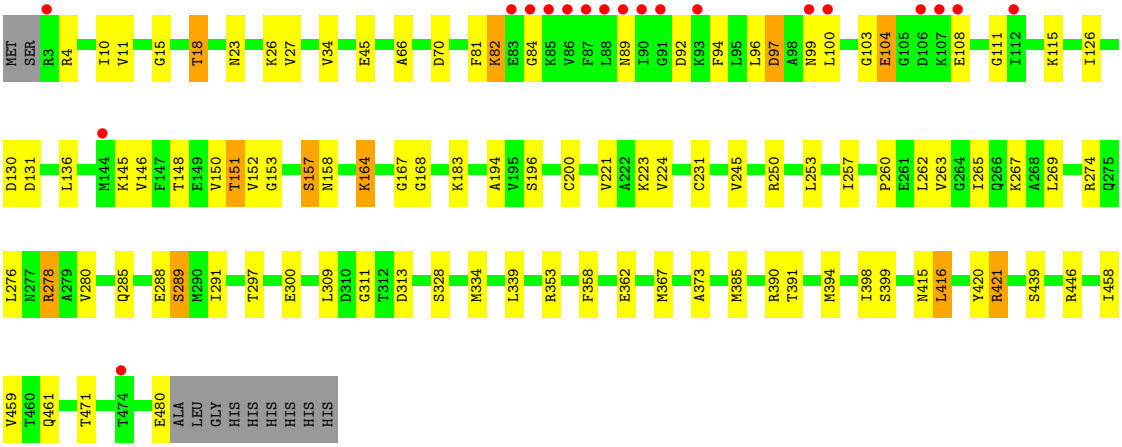
• Molecule 1: Pyruvate kinase

Chain A:



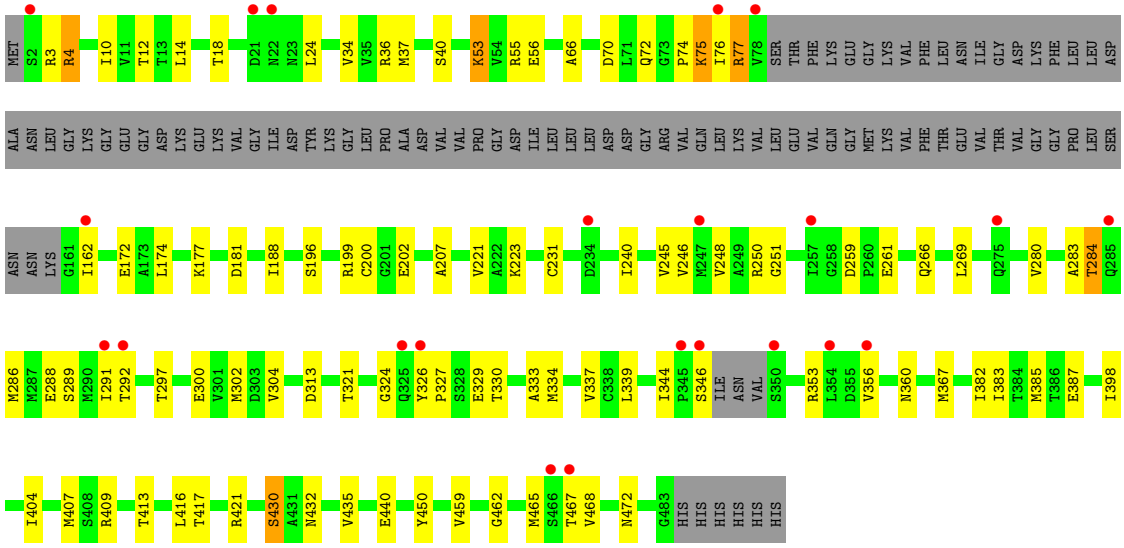
• Molecule 1: Pyruvate kinase

Chain B:



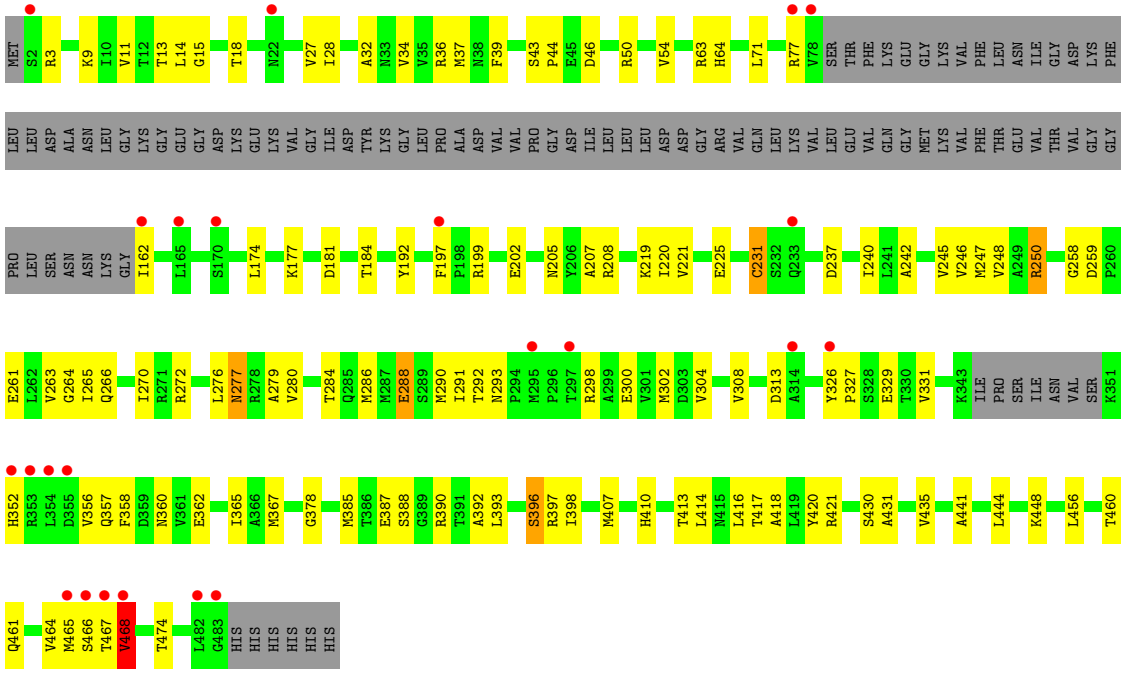
• Molecule 1: Pyruvate kinase

Chain C:



- Molecule 1: Pyruvate kinase

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	117.69Å 137.30Å 139.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.95 – 2.68 32.95 – 2.68	Depositor EDS
% Data completeness (in resolution range)	98.5 (32.95-2.68) 98.6 (32.95-2.68)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.51 (at 2.68Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.210 , 0.259 0.210 , 0.259	Depositor DCC
R_{free} test set	3101 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å^2)	51.5	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 46.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.017 for -h,l,k	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	13318	wwPDB-VP
Average B, all atoms (Å^2)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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: MG, CIT

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	0.47	0/3615	0.68	2/4882 (0.0%)
1	B	0.45	0/3609	0.66	1/4874 (0.0%)
1	C	0.42	0/2981	0.64	2/4025 (0.0%)
1	D	0.40	0/2949	0.59	0/3981
All	All	0.44	0/13154	0.65	5/17762 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	156	LEU	CB-CG-CD2	-6.71	99.60	111.00
1	C	416	LEU	CA-CB-CG	5.59	128.16	115.30
1	B	416	LEU	CA-CB-CG	5.45	127.83	115.30
1	A	95	LEU	CA-CB-CG	5.36	127.63	115.30
1	C	416	LEU	CB-CG-CD1	-5.18	102.19	111.00

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	3579	0	3678	82	0
1	B	3573	0	3674	66	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2955	0	3029	55	0
1	D	2924	0	2998	70	0
2	A	1	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0
3	A	13	0	5	0	0
3	C	13	0	5	1	0
3	D	13	0	5	1	0
4	A	77	0	0	3	0
4	B	105	0	0	4	0
4	C	47	0	0	1	0
4	D	15	0	0	1	0
All	All	13318	0	13394	267	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 10.

All (267) 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:81:PHE:HB3	1:A:104:GLU:HA	1.49	0.94
1:B:82:LYS:HE2	1:B:104:GLU:HB2	1.60	0.84
1:B:100:LEU:HD21	1:B:103:GLY:H	1.44	0.81
1:D:231:CYS:SG	4:D:612:HOH:O	2.39	0.80
1:A:107:LYS:HD2	1:A:109:LYS:HE3	1.64	0.79
1:C:291:ILE:HD12	1:C:324:GLY:HA2	1.66	0.78
1:B:157:SER:OG	1:B:158:ASN:N	2.14	0.78
1:C:288:GLU:O	1:C:291:ILE:HG22	1.85	0.77
1:C:250:ARG:NH1	1:D:302:MET:SD	2.58	0.76
1:A:135:GLN:HB2	1:A:152:VAL:HG12	1.67	0.76
1:A:95:LEU:HB3	1:A:108:GLU:O	1.88	0.74
1:B:221:VAL:HG22	1:B:245:VAL:HB	1.71	0.73
1:A:231:CYS:SG	4:A:672:HOH:O	2.47	0.73
1:B:34:VAL:HG22	1:B:66:ALA:HB3	1.70	0.73
1:C:77:ARG:HH11	1:C:77:ARG:HG2	1.55	0.71
1:D:208:ARG:NH2	1:D:242:ALA:O	2.23	0.71
1:C:14:LEU:HD22	1:C:18:THR:HG21	1.72	0.70
1:A:200:CYS:SG	4:A:675:HOH:O	2.48	0.69
1:A:326:TYR:HB3	1:A:329:GLU:HB2	1.75	0.69
1:B:399:SER:HA	1:B:421:ARG:NH2	2.07	0.69
1:B:183:LYS:NZ	4:B:502:HOH:O	2.25	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:GLY:HA2	1:A:109:LYS:O	1.94	0.67
1:D:392:ALA:O	1:D:396:SER:OG	2.12	0.67
1:D:460:THR:HG22	1:D:474:THR:HG22	1.74	0.67
1:A:297:THR:HG22	1:A:300:GLU:H	1.60	0.67
1:B:399:SER:HA	1:B:421:ARG:HH21	1.59	0.67
1:B:200:CYS:SG	4:B:603:HOH:O	2.51	0.66
1:A:98:ALA:HB3	1:A:144:MET:HB3	1.79	0.64
1:B:11:VAL:HG22	1:B:34:VAL:HB	1.79	0.64
1:C:353:ARG:HD3	1:C:356:VAL:HG11	1.80	0.64
1:C:34:VAL:HG22	1:C:66:ALA:HB3	1.80	0.64
1:A:105:GLY:O	1:A:107:LYS:N	2.31	0.64
1:A:96:LEU:HB2	1:A:146:VAL:HG13	1.79	0.64
1:A:407:MET:HE1	1:A:445:LEU:HD11	1.79	0.64
1:B:390:ARG:O	1:B:394:MET:HG3	1.98	0.63
1:A:4:ARG:NH2	1:A:339:LEU:O	2.31	0.63
1:A:81:PHE:HD1	1:A:105:GLY:H	1.47	0.63
1:D:39:PHE:HD2	1:D:71:LEU:HD23	1.64	0.63
1:C:387:GLU:O	1:C:413:THR:HG21	2.00	0.62
1:D:326:TYR:HB3	1:D:329:GLU:HB2	1.79	0.62
1:C:76:ILE:HG22	1:C:162:ILE:HG13	1.81	0.62
1:D:219:LYS:HE2	1:D:416:LEU:HD21	1.81	0.62
1:C:291:ILE:HG23	1:C:292:THR:HG23	1.80	0.62
1:C:313:ASP:HA	1:C:421:ARG:HB2	1.80	0.62
1:B:164:LYS:HE3	1:B:168:GLY:HA3	1.82	0.61
1:C:248:VAL:HG21	1:C:266:GLN:HG2	1.82	0.61
1:C:12:THR:HB	1:C:321:THR:HG21	1.81	0.61
1:C:367:MET:SD	1:C:398:ILE:HD11	2.41	0.61
1:D:9:LYS:NZ	1:D:420:TYR:O	2.34	0.61
1:D:464:VAL:O	1:D:466:SER:N	2.33	0.61
1:C:283:ALA:O	1:C:284:THR:HG22	2.01	0.60
1:B:313:ASP:HA	1:B:421:ARG:HB2	1.84	0.60
1:B:353:ARG:HD2	1:B:367:MET:HE2	1.83	0.60
1:D:248:VAL:HG21	1:D:266:GLN:HG2	1.83	0.60
1:B:362:GLU:HG3	1:B:471:THR:HB	1.83	0.59
1:A:81:PHE:HD2	1:A:158:ASN:ND2	2.00	0.59
1:A:23:ASN:O	1:A:27:VAL:HG13	2.03	0.59
1:A:130:ASP:O	1:A:133:ARG:HG3	2.02	0.59
1:B:4:ARG:NH2	1:B:339:LEU:O	2.35	0.59
1:C:467:THR:HG22	1:C:468:VAL:H	1.68	0.57
1:D:387:GLU:HG3	1:D:410:HIS:CD2	2.39	0.57
1:D:413:THR:O	1:D:417:THR:HG23	2.04	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:358:PHE:CZ	1:A:367:MET:HG3	2.40	0.57
1:B:89:ASN:HB2	1:B:150:VAL:HB	1.85	0.57
1:C:174:LEU:HD11	1:C:207:ALA:HB2	1.86	0.57
1:A:81:PHE:CB	1:A:104:GLU:HA	2.30	0.57
1:B:285:GLN:HE21	1:B:288:GLU:HG3	1.68	0.57
1:B:23:ASN:O	1:B:27:VAL:HG13	2.05	0.57
1:D:28:ILE:O	1:D:63:ARG:NH1	2.37	0.57
1:B:297:THR:HG22	1:B:300:GLU:H	1.70	0.56
1:C:4:ARG:NH2	1:C:339:LEU:O	2.38	0.56
1:C:74:PRO:HG3	1:C:172:GLU:O	2.05	0.56
1:B:96:LEU:HB2	1:B:146:VAL:HG13	1.88	0.55
1:C:24:LEU:HD23	1:C:53:LYS:HD2	1.88	0.55
1:A:81:PHE:CD2	1:A:158:ASN:ND2	2.74	0.55
1:A:87:PHE:CE1	1:A:150:VAL:HG22	2.42	0.55
1:A:131:ASP:HA	1:A:255:VAL:HG21	1.88	0.55
1:A:139:LEU:HD21	1:A:149:GLU:HB3	1.87	0.55
1:B:96:LEU:HB2	1:B:146:VAL:CG1	2.36	0.55
1:C:432:ASN:ND2	1:C:440:GLU:OE2	2.36	0.55
1:B:250:ARG:HB3	1:B:262:LEU:HD21	1.89	0.55
1:A:97:ASP:O	1:A:111:GLY:HA2	2.07	0.54
1:B:97:ASP:HB2	1:B:145:LYS:HG2	1.89	0.54
1:B:26:LYS:HB3	1:B:328:SER:HB3	1.88	0.54
1:B:288:GLU:OE2	4:B:501:HOH:O	2.18	0.54
1:A:251:GLY:H	1:A:284:THR:HG21	1.72	0.54
1:B:94:PHE:CD1	1:B:108:GLU:HA	2.42	0.54
1:D:64:HIS:NE2	1:D:414:LEU:HD13	2.22	0.54
1:B:289:SER:HB3	1:B:300:GLU:OE1	2.07	0.54
1:D:177:LYS:NZ	1:D:181:ASP:OD2	2.39	0.54
1:B:385:MET:HE3	1:B:461:GLN:HB3	1.90	0.53
1:D:50:ARG:O	1:D:54:VAL:HG23	2.07	0.53
1:A:50:ARG:HA	1:A:53:LYS:HG2	1.90	0.53
1:A:287:MET:HG3	1:A:334:MET:CE	2.39	0.53
1:A:107:LYS:HG3	1:A:109:LYS:HG3	1.91	0.52
1:C:333:ALA:O	1:C:337:VAL:HG23	2.10	0.52
1:A:267:LYS:HE2	1:B:309:LEU:HD13	1.91	0.52
1:B:94:PHE:HA	1:B:108:GLU:HB2	1.91	0.52
1:C:37:MET:HE1	1:C:188:ILE:HD13	1.91	0.52
1:A:81:PHE:HD1	1:A:105:GLY:N	2.07	0.52
1:A:430:SER:OG	1:A:440:GLU:OE1	2.25	0.52
1:C:326:TYR:HB3	1:C:329:GLU:HB2	1.92	0.52
1:A:250:ARG:H	1:A:284:THR:HG23	1.74	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:353:ARG:HD2	1:B:367:MET:CE	2.40	0.52
1:C:10:ILE:HG12	1:C:334:MET:HE3	1.92	0.51
1:B:151:THR:OG1	1:B:152:VAL:N	2.41	0.51
1:B:224:VAL:HG11	1:B:269:LEU:HD11	1.92	0.51
1:C:344:ILE:HG21	1:D:264:GLY:HA3	1.92	0.51
1:D:367:MET:CE	1:D:398:ILE:HD11	2.39	0.51
1:A:284:THR:OG1	1:A:285:GLN:N	2.44	0.51
1:B:89:ASN:HD21	1:B:153:GLY:N	2.08	0.51
1:A:96:LEU:HB2	1:A:146:VAL:CG1	2.41	0.51
1:D:270:ILE:HG12	1:D:280:VAL:HG11	1.91	0.51
1:A:267:LYS:HE2	1:B:309:LEU:CD1	2.41	0.51
1:A:240:ILE:HG12	1:A:246:VAL:HG21	1.92	0.51
1:A:151:THR:OG1	1:A:152:VAL:N	2.39	0.50
3:D:502:CIT:O3	3:D:502:CIT:O7	2.26	0.50
1:D:358:PHE:CE1	1:D:367:MET:HG3	2.46	0.50
1:C:200:CYS:SG	1:C:202:GLU:HB3	2.52	0.50
1:A:123:PRO:HB3	1:A:139:LEU:O	2.11	0.50
1:C:261:GLU:OE2	1:C:261:GLU:N	2.37	0.49
1:D:14:LEU:HD12	1:D:37:MET:HE1	1.94	0.49
1:C:196:SER:HA	1:C:223:LYS:HD3	1.94	0.49
1:A:82:LYS:H	1:A:104:GLU:HB3	1.76	0.49
1:A:140:GLU:HG3	1:A:147:PHE:CD1	2.47	0.49
1:B:276:LEU:O	1:B:278:ARG:HD3	2.12	0.49
1:D:467:THR:O	1:D:468:VAL:HG12	2.13	0.49
1:D:13:THR:HA	1:D:36:ARG:HB3	1.95	0.49
1:A:208:ARG:CZ	1:A:220:ILE:HD13	2.43	0.49
1:A:79:SER:OG	1:A:80:THR:N	2.46	0.48
1:B:263:VAL:HG12	1:B:267:LYS:HE3	1.95	0.48
1:C:462:GLY:HA2	1:C:472:ASN:OD1	2.13	0.48
1:D:288:GLU:HG3	1:D:291:ILE:HD12	1.95	0.48
1:A:83:GLU:HA	1:A:85:LYS:HZ3	1.77	0.48
1:A:81:PHE:CD2	1:A:81:PHE:N	2.81	0.48
1:C:231:CYS:SG	4:C:633:HOH:O	2.61	0.48
1:A:251:GLY:O	1:A:255:VAL:HG23	2.14	0.48
1:B:373:ALA:HB2	1:B:458:ILE:HG13	1.95	0.48
1:C:297:THR:HG23	1:C:300:GLU:H	1.79	0.48
1:D:245:VAL:HG22	1:D:279:ALA:HB3	1.95	0.48
1:A:87:PHE:HE1	1:A:150:VAL:HG13	1.78	0.47
1:B:367:MET:CE	1:B:398:ILE:HD11	2.44	0.47
1:D:221:VAL:CG1	1:D:247:MET:HB2	2.44	0.47
1:D:430:SER:O	1:D:431:ALA:HB3	2.14	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:PHE:HD2	1:A:158:ASN:HD22	1.60	0.47
1:A:81:PHE:H	1:A:158:ASN:HD21	1.61	0.47
1:B:385:MET:HE2	1:B:459:VAL:HG11	1.96	0.47
1:C:385:MET:HE2	1:C:459:VAL:HG11	1.95	0.47
1:A:414:LEU:HD23	4:A:651:HOH:O	2.14	0.47
1:D:237:ASP:OD2	1:D:272:ARG:NH2	2.47	0.47
1:A:430:SER:O	1:A:431:ALA:HB3	2.14	0.47
1:C:251:GLY:HA2	1:D:298:ARG:HH22	1.80	0.47
1:C:286:MET:HE3	1:C:304:VAL:HG22	1.96	0.47
1:B:89:ASN:HD21	1:B:152:VAL:C	2.18	0.47
1:B:196:SER:OG	1:B:223:LYS:NZ	2.41	0.47
1:D:304:VAL:O	1:D:308:VAL:HG23	2.15	0.47
1:C:344:ILE:HG22	1:C:346:SER:H	1.78	0.47
1:C:3:ARG:HD2	1:C:4:ARG:H	1.80	0.47
1:C:37:MET:CE	1:C:188:ILE:HD13	2.44	0.47
1:D:9:LYS:HG3	1:D:418:ALA:O	2.14	0.47
1:D:313:ASP:HA	1:D:421:ARG:HB2	1.96	0.46
1:D:385:MET:HE3	1:D:461:GLN:CB	2.45	0.46
1:A:251:GLY:N	1:A:284:THR:HG21	2.30	0.46
1:D:221:VAL:HG12	1:D:247:MET:HB2	1.96	0.46
1:B:10:ILE:HG12	1:B:334:MET:HG2	1.98	0.46
1:D:15:GLY:O	1:D:18:THR:HB	2.15	0.46
1:B:126:ILE:HA	1:B:136:LEU:O	2.16	0.46
1:B:92:ASP:O	1:B:150:VAL:HG23	2.16	0.45
1:C:221:VAL:HG22	1:C:245:VAL:HB	1.98	0.45
1:D:39:PHE:HB3	1:D:181:ASP:OD2	2.16	0.45
1:B:66:ALA:HB2	1:B:415:ASN:O	2.17	0.45
1:A:87:PHE:CE1	1:A:150:VAL:HG13	2.51	0.45
1:A:253:LEU:O	1:A:257:ILE:HG12	2.17	0.45
1:B:274:ARG:HD2	1:B:311:GLY:O	2.17	0.45
1:B:285:GLN:NE2	1:B:288:GLU:HG3	2.30	0.45
1:A:274:ARG:HD2	1:A:311:GLY:O	2.16	0.45
1:C:177:LYS:NZ	1:C:181:ASP:OD2	2.44	0.45
1:D:407:MET:CE	1:D:441:ALA:HB1	2.46	0.45
1:A:353:ARG:HD2	1:A:367:MET:SD	2.56	0.45
1:B:94:PHE:O	1:B:148:THR:HG22	2.17	0.45
1:C:289:SER:OG	1:C:300:GLU:OE1	2.30	0.45
1:D:11:VAL:HG22	1:D:34:VAL:HB	1.99	0.45
1:C:75:LYS:HD2	1:C:77:ARG:HD3	1.98	0.45
1:D:44:PRO:HG3	1:D:184:THR:HG21	1.99	0.45
1:C:326:TYR:O	1:C:330:THR:OG1	2.26	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:277:ASN:HA	1:D:397:ARG:HG3	1.99	0.44
1:C:302:MET:HG2	1:D:263:VAL:HG22	1.99	0.44
1:C:385:MET:HE2	1:C:459:VAL:CG1	2.47	0.44
1:D:27:VAL:HB	1:D:331:VAL:HG21	1.99	0.44
1:C:450:TYR:OH	3:C:503:CIT:O5	2.18	0.44
1:D:248:VAL:CG2	1:D:266:GLN:HG2	2.48	0.44
1:A:83:GLU:HA	1:A:85:LYS:NZ	2.32	0.44
1:B:446:ARG:HH11	1:B:480:GLU:HB2	1.82	0.44
1:D:32:ALA:O	1:D:63:ARG:NH2	2.51	0.44
1:D:208:ARG:CZ	1:D:220:ILE:HD12	2.48	0.44
1:A:39:PHE:O	1:A:177:LYS:NZ	2.48	0.44
1:B:97:ASP:O	1:B:111:GLY:HA2	2.18	0.44
1:A:241:LEU:HA	1:A:278:ARG:NH2	2.32	0.43
1:B:70:ASP:HA	1:B:194:ALA:HB3	2.00	0.43
1:B:367:MET:HE3	1:B:398:ILE:HD11	2.00	0.43
1:C:248:VAL:HB	1:C:269:LEU:HD23	1.99	0.43
1:A:15:GLY:H	1:A:18:THR:HG22	1.82	0.43
1:B:280:VAL:O	1:B:313:ASP:HB2	2.17	0.43
1:B:358:PHE:HA	4:B:552:HOH:O	2.18	0.43
1:D:13:THR:OG1	1:D:36:ARG:NH1	2.50	0.43
1:A:24:LEU:O	1:A:27:VAL:HG22	2.19	0.43
1:C:430:SER:HB2	1:C:440:GLU:OE1	2.18	0.43
1:D:3:ARG:HA	1:D:3:ARG:HD2	1.65	0.43
1:D:192:TYR:OH	1:D:416:LEU:HD12	2.19	0.43
1:D:290:MET:HA	1:D:293:ASN:O	2.18	0.43
1:A:458:ILE:HD13	1:A:476:ILE:HG12	2.01	0.43
1:C:40:SER:OG	1:C:72:GLN:HG3	2.18	0.43
1:D:393:LEU:O	1:D:397:ARG:HG2	2.18	0.43
1:D:387:GLU:HG3	1:D:410:HIS:NE2	2.33	0.43
1:D:388:SER:OG	1:D:390:ARG:HG3	2.19	0.43
1:D:258:GLY:HA3	1:D:261:GLU:OE2	2.18	0.43
1:A:37:MET:HE1	1:A:67:ILE:HD13	2.01	0.43
1:A:93:LYS:O	1:A:108:GLU:HB3	2.18	0.43
1:A:340:GLY:HA3	1:B:260:PRO:HB3	2.00	0.43
1:A:365:ILE:HG13	1:A:472:ASN:HA	2.01	0.43
1:B:82:LYS:HG3	1:B:104:GLU:H	1.82	0.43
1:C:382:ILE:O	1:C:404:ILE:HA	2.18	0.43
1:D:162:ILE:HD12	1:D:162:ILE:HA	1.89	0.43
1:B:253:LEU:O	1:B:257:ILE:HG12	2.19	0.42
1:C:383:ILE:HB	1:C:459:VAL:HG22	2.00	0.42
1:D:444:LEU:HD11	1:D:448:LYS:HE2	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:MET:HE2	1:A:330:THR:HG23	2.02	0.42
1:B:99:ASN:O	1:B:100:LEU:HB2	2.18	0.42
1:A:221:VAL:HG22	1:A:245:VAL:HB	2.01	0.42
1:A:251:GLY:H	1:A:284:THR:CG2	2.32	0.42
1:C:284:THR:HG23	1:C:286:MET:HG3	2.01	0.42
1:D:365:ILE:HG23	1:D:474:THR:HG23	2.01	0.42
1:B:15:GLY:O	1:B:18:THR:HG22	2.20	0.42
1:B:416:LEU:HD21	1:B:420:TYR:OH	2.20	0.42
1:A:51:ALA:O	1:A:55:ARG:HG3	2.20	0.42
1:D:202:GLU:HG2	1:D:205:ASN:ND2	2.35	0.42
1:A:8:THR:HG21	1:A:308:VAL:HG22	2.02	0.42
1:A:208:ARG:HG3	1:A:220:ILE:HD11	2.02	0.42
1:A:2:SER:N	1:A:449:GLY:O	2.53	0.41
1:C:240:ILE:HG12	1:C:246:VAL:HG21	2.01	0.41
1:B:89:ASN:CB	1:B:150:VAL:HB	2.49	0.41
1:B:131:ASP:N	1:B:131:ASP:OD1	2.53	0.41
1:D:378:GLY:O	1:D:456:LEU:HB2	2.21	0.41
1:A:79:SER:HB2	1:A:113:ASP:OD1	2.19	0.41
1:B:285:GLN:HE22	1:B:291:ILE:HD11	1.85	0.41
1:C:413:THR:O	1:C:417:THR:HG23	2.20	0.41
1:D:240:ILE:HG12	1:D:246:VAL:HG21	2.01	0.41
1:A:87:PHE:HD1	1:A:88:LEU:O	2.03	0.41
1:A:313:ASP:HA	1:A:421:ARG:HB2	2.02	0.41
1:A:459:VAL:O	1:A:474:THR:HA	2.20	0.41
1:C:36:ARG:NE	1:C:70:ASP:OD2	2.41	0.41
1:D:174:LEU:HD11	1:D:207:ALA:HB2	2.02	0.41
1:D:43:SER:O	1:D:46:ASP:N	2.51	0.41
1:D:276:LEU:HA	1:D:276:LEU:HD23	1.86	0.41
1:A:172:GLU:HA	1:A:172:GLU:OE1	2.21	0.41
1:B:265:ILE:HD13	1:B:265:ILE:HA	1.87	0.41
1:D:39:PHE:CD2	1:D:71:LEU:HD23	2.49	0.41
1:A:117:LEU:N	1:A:118:PRO:HD2	2.36	0.41
1:D:286:MET:O	1:D:300:GLU:HG2	2.21	0.41
1:D:356:VAL:HG22	1:D:357:GLN:N	2.35	0.41
1:A:412:ARG:O	1:A:416:LEU:HB2	2.21	0.40
1:A:353:ARG:HB3	1:A:356:VAL:HB	2.03	0.40
1:D:77:ARG:HH11	1:D:77:ARG:HD3	1.68	0.40
1:A:228:GLU:OE2	1:A:228:GLU:N	2.53	0.40
1:D:250:ARG:HB2	1:D:284:THR:OG1	2.21	0.40
1:D:261:GLU:O	1:D:265:ILE:HG12	2.21	0.40
1:D:360:ASN:ND2	1:D:362:GLU:HB2	2.36	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	477/489 (98%)	437 (92%)	33 (7%)	7 (2%)	10	23
1	B	476/489 (97%)	438 (92%)	32 (7%)	6 (1%)	12	27
1	C	391/489 (80%)	368 (94%)	22 (6%)	1 (0%)	41	64
1	D	386/489 (79%)	362 (94%)	20 (5%)	4 (1%)	15	34
All	All	1730/1956 (88%)	1605 (93%)	107 (6%)	18 (1%)	15	34

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	87	PHE
1	A	106	ASP
1	A	107	LYS
1	B	104	GLU
1	D	352	HIS
1	D	465	MET
1	D	468	VAL
1	A	284	THR
1	B	82	LYS
1	B	167	GLY
1	A	90	ILE
1	B	84	GLY
1	A	83	GLU
1	D	327	PRO
1	A	105	GLY
1	B	130	ASP
1	B	151	THR
1	C	327	PRO

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	383/391 (98%)	371 (97%)	12 (3%)	40	67
1	B	382/391 (98%)	369 (97%)	13 (3%)	37	63
1	C	313/391 (80%)	297 (95%)	16 (5%)	24	46
1	D	309/391 (79%)	297 (96%)	12 (4%)	32	58
All	All	1387/1564 (89%)	1334 (96%)	53 (4%)	33	59

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ARG
1	A	18	THR
1	A	92	ASP
1	A	107	LYS
1	A	158	ASN
1	A	172	GLU
1	A	225	GLU
1	A	334	MET
1	A	386	THR
1	A	394	MET
1	A	435	VAL
1	A	465	MET
1	B	18	THR
1	B	45	GLU
1	B	81	PHE
1	B	97	ASP
1	B	115	LYS
1	B	157	SER
1	B	164	LYS
1	B	231	CYS
1	B	278	ARG
1	B	289	SER
1	B	391	THR
1	B	421	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	439	SER
1	C	4	ARG
1	C	53	LYS
1	C	55	ARG
1	C	56	GLU
1	C	75	LYS
1	C	77	ARG
1	C	199	ARG
1	C	259	ASP
1	C	280	VAL
1	C	284	THR
1	C	360	ASN
1	C	407	MET
1	C	409	ARG
1	C	430	SER
1	C	435	VAL
1	C	465	MET
1	D	197	PHE
1	D	199	ARG
1	D	225	GLU
1	D	231	CYS
1	D	250	ARG
1	D	259	ASP
1	D	277	ASN
1	D	288	GLU
1	D	292	THR
1	D	396	SER
1	D	435	VAL
1	D	468	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	B	89	ASN
1	B	285	GLN
1	D	266	GLN
1	D	277	ASN
1	D	360	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 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	CIT	A	502	-	12,12,12	1.57	2 (16%)	17,17,17	2.38	10 (58%)
3	CIT	C	503	-	12,12,12	1.46	1 (8%)	17,17,17	2.26	7 (41%)
3	CIT	D	502	-	12,12,12	1.01	0	17,17,17	1.15	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CIT	A	502	-	-	6/16/16/16	-
3	CIT	C	503	-	-	5/16/16/16	-
3	CIT	D	502	-	-	12/16/16/16	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	CIT	C3-C6	3.56	1.57	1.53
3	C	503	CIT	C3-C6	3.24	1.56	1.53
3	A	502	CIT	O7-C3	2.12	1.47	1.43

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	503	CIT	O6-C6-C3	5.18	122.05	113.05
3	A	502	CIT	O7-C3-C6	-3.63	103.77	108.86
3	A	502	CIT	O5-C6-C3	3.53	127.25	122.25
3	A	502	CIT	O2-C1-O1	-3.38	114.86	123.30
3	C	503	CIT	O7-C3-C4	-3.22	101.88	109.40
3	A	502	CIT	O2-C1-C2	3.17	124.55	114.35
3	C	503	CIT	O2-C1-O1	-3.15	115.44	123.30
3	C	503	CIT	O6-C6-O5	-3.09	113.98	123.82
3	A	502	CIT	O6-C6-O5	-2.64	115.43	123.82
3	A	502	CIT	C3-C4-C5	-2.49	107.78	113.81
3	A	502	CIT	O6-C6-C3	2.46	117.31	113.05
3	A	502	CIT	O4-C5-C4	2.44	122.19	114.35
3	A	502	CIT	C4-C3-C2	-2.37	102.97	109.16
3	C	503	CIT	C3-C4-C5	2.25	119.25	113.81
3	C	503	CIT	O4-C5-O3	-2.20	117.82	123.30
3	A	502	CIT	O7-C3-C2	2.15	114.44	109.40
3	C	503	CIT	O7-C3-C2	-2.14	104.38	109.40

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	CIT	C1-C2-C3-O7
3	A	502	CIT	C1-C2-C3-C4
3	A	502	CIT	C1-C2-C3-C6
3	C	503	CIT	C1-C2-C3-O7
3	C	503	CIT	C1-C2-C3-C4
3	C	503	CIT	C1-C2-C3-C6
3	D	502	CIT	C2-C3-C4-C5
3	D	502	CIT	O7-C3-C4-C5
3	D	502	CIT	C6-C3-C4-C5
3	D	502	CIT	O7-C3-C6-O5
3	D	502	CIT	O7-C3-C6-O6
3	D	502	CIT	C4-C3-C6-O5
3	D	502	CIT	C4-C3-C6-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	D	502	CIT	C1-C2-C3-O7
3	D	502	CIT	C1-C2-C3-C4
3	D	502	CIT	C1-C2-C3-C6
3	A	502	CIT	C6-C3-C4-C5
3	C	503	CIT	C2-C3-C4-C5
3	C	503	CIT	O7-C3-C4-C5
3	D	502	CIT	C2-C3-C6-O6
3	D	502	CIT	C2-C3-C6-O5
3	A	502	CIT	O7-C3-C4-C5
3	A	502	CIT	C2-C3-C4-C5

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	503	CIT	1	0
3	D	502	CIT	1	0

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	479/489 (97%)	0.16	32 (6%)	17	15	29, 45, 97, 148	0
1	B	478/489 (97%)	0.06	19 (3%)	38	36	30, 44, 106, 152	0
1	C	397/489 (81%)	0.19	22 (5%)	25	23	36, 57, 89, 123	0
1	D	392/489 (80%)	0.29	23 (5%)	22	20	38, 67, 93, 131	0
All	All	1746/1956 (89%)	0.17	96 (5%)	25	23	29, 53, 97, 152	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	84	GLY	8.3
1	B	90	ILE	8.0
1	B	108	GLU	7.9
1	C	466	SER	6.6
1	A	106	ASP	6.5
1	A	104	GLU	6.3
1	A	86	VAL	5.9
1	D	78	VAL	5.8
1	C	350	SER	5.6
1	A	88	LEU	5.6
1	B	84	GLY	4.9
1	A	87	PHE	4.7
1	B	87	PHE	4.6
1	B	88	LEU	4.6
1	B	89	ASN	4.5
1	A	89	ASN	4.2
1	C	21	ASP	4.1
1	C	162	ILE	4.0
1	A	108	GLU	3.9
1	A	83	GLU	3.9
1	C	2	SER	3.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	77	ARG	3.8
1	D	468	VAL	3.7
1	A	80	THR	3.6
1	B	91	GLY	3.6
1	C	326	TYR	3.5
1	A	107	LYS	3.5
1	C	345	PRO	3.5
1	A	81	PHE	3.5
1	A	151	THR	3.4
1	A	155	PRO	3.4
1	C	346	SER	3.2
1	C	467	THR	3.2
1	A	144	MET	3.2
1	C	292	THR	3.2
1	B	144	MET	3.1
1	D	354	LEU	3.1
1	C	325	GLN	3.0
1	D	353	ARG	3.0
1	D	295	MET	3.0
1	B	100	LEU	3.0
1	B	106	ASP	3.0
1	B	83	GLU	3.0
1	A	103	GLY	2.9
1	D	352	HIS	2.9
1	B	3	ARG	2.9
1	A	101	GLY	2.8
1	B	93	LYS	2.8
1	B	99	ASN	2.8
1	A	142	GLN	2.7
1	B	85	LYS	2.7
1	D	165	LEU	2.7
1	B	86	VAL	2.6
1	B	107	LYS	2.6
1	C	275	GLN	2.6
1	A	141	VAL	2.5
1	A	115	LYS	2.5
1	B	112	ILE	2.5
1	A	21	ASP	2.5
1	C	354	LEU	2.5
1	D	22	ASN	2.5
1	A	284	THR	2.4
1	D	466	SER	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	285	GLN	2.4
1	D	314	ALA	2.4
1	D	2	SER	2.4
1	C	234	ASP	2.4
1	C	76	ILE	2.3
1	C	257	ILE	2.3
1	C	22	ASN	2.3
1	C	356	VAL	2.3
1	D	467	THR	2.3
1	D	465	MET	2.3
1	C	247	MET	2.2
1	D	355	ASP	2.2
1	A	85	LYS	2.2
1	A	370	MET	2.2
1	A	452	MET	2.2
1	A	3	ARG	2.2
1	D	297	THR	2.2
1	D	482	LEU	2.2
1	D	483	GLY	2.2
1	A	90	ILE	2.2
1	A	100	LEU	2.1
1	D	197	PHE	2.1
1	D	233	GLN	2.1
1	A	105	GLY	2.1
1	A	82	LYS	2.1
1	A	145	LYS	2.1
1	B	474	THR	2.1
1	D	326	TYR	2.1
1	C	78	VAL	2.1
1	A	292	THR	2.0
1	D	170	SER	2.0
1	C	291	ILE	2.0
1	D	162	ILE	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
3	CIT	A	502	13/13	0.88	0.16	43,50,60,60	0
3	CIT	D	502	13/13	0.88	0.14	57,66,71,74	0
3	CIT	C	503	13/13	0.92	0.14	44,54,59,66	0
2	MG	C	501	1/1	0.92	0.17	58,58,58,58	0
2	MG	D	501	1/1	0.94	0.26	51,51,51,51	0
2	MG	C	502	1/1	0.94	0.40	31,31,31,31	0
2	MG	A	501	1/1	0.98	0.13	43,43,43,43	0

6.5 Other polymers (i)

There are no such residues in this entry.