Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 11:00 PM EDT

PDB ID	:	1EJX
Title	:	CRYSTAL STRUCTURE OF WILD-TYPE KLEBSIELLA AEROGENES UREASE AT 100K
Authors	:	Pearson, M.A.; Karplus, P.A.
Deposited on	:	2000-03-04
Resolution	:	1.60 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:

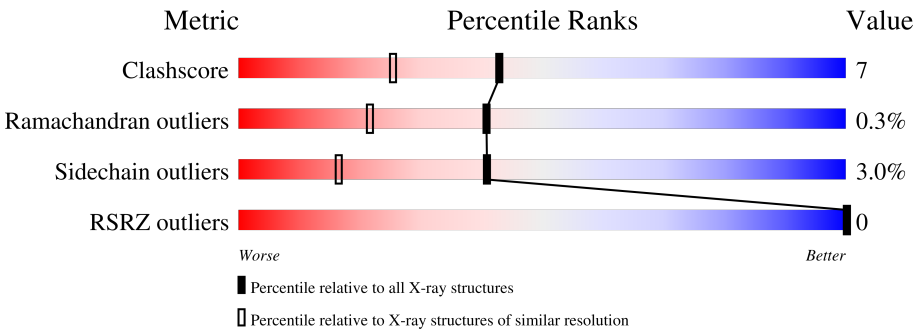
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 1.60 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))
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	C	567	82% 14% • •
2	B	101	72% 25% •
3	A	100	80% 19% •

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 6272 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 UREASE ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	C	556	4161	2615	727	796	23	0	5	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1217	KCX	LYS	MODIFIED RESIDUE	UNP P18314

- Molecule 2 is a protein called UREASE BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	101	785	496	150	136	3	0	0	0

- Molecule 3 is a protein called UREASE GAMMA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	A	100	776	491	134	146	5	0	0	0

- Molecule 4 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ni		
4	C	2	2	2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	C	402	402	402	0	0

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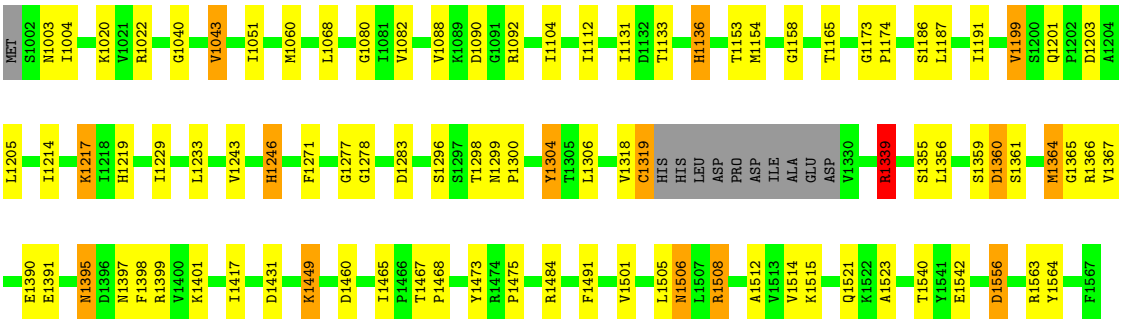
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	B	76	76	76	0	0
			Total	O		
5	A	70	70	70	0	0

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.

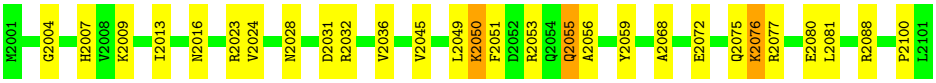
- Molecule 1: UREASE ALPHA SUBUNIT

Chain C:



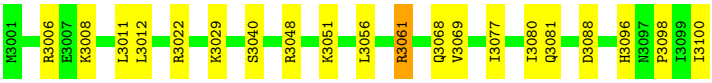
- Molecule 2: UREASE BETA SUBUNIT

Chain B:



- Molecule 3: UREASE GAMMA SUBUNIT

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α, β, γ	168.96Å 168.96Å 168.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.60 39.82 – 1.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-1.60) 92.8 (39.82-1.60)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.19 (at 1.60Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.172 , 0.240 0.171 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	18.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 105.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.028 for -l,-k,-h	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	6272	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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: KCX, NI

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	C	0.98	1/4254 (0.0%)	1.38	32/5794 (0.6%)
2	B	0.96	0/805	1.42	6/1087 (0.6%)
3	A	1.04	0/787	1.37	7/1061 (0.7%)
All	All	0.99	1/5846 (0.0%)	1.38	45/7942 (0.6%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1556	ASP	CB-CG	6.40	1.65	1.51

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	1092	ARG	NE-CZ-NH2	-18.30	111.15	120.30
1	C	1092	ARG	NE-CZ-NH1	12.82	126.71	120.30
1	C	1092	ARG	CD-NE-CZ	10.73	138.62	123.60
1	C	1556	ASP	CB-CG-OD1	8.81	126.23	118.30
1	C	1556	ASP	N-CA-CB	-8.13	95.97	110.60
2	B	2088	ARG	NE-CZ-NH2	-7.07	116.77	120.30
3	A	3068	GLN	N-CA-CB	-7.04	97.92	110.60
1	C	1484	ARG	NE-CZ-NH1	6.91	123.76	120.30
2	B	2056	ALA	N-CA-CB	-6.83	100.54	110.10
1	C	1339	ARG	NE-CZ-NH1	6.73	123.67	120.30
1	C	1564	TYR	CG-CD1-CE1	-6.72	115.92	121.30
1	C	1092	ARG	CG-CD-NE	-6.41	98.33	111.80
1	C	1299	ASN	N-CA-C	6.36	128.16	111.00
1	C	1460	ASP	CB-CG-OD1	6.25	123.93	118.30
2	B	2032	ARG	NE-CZ-NH1	6.25	123.42	120.30
3	A	3040	SER	CB-CA-C	6.12	121.73	110.10
2	B	2100	PRO	C-N-CA	6.10	136.95	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	1473	TYR	CB-CG-CD2	-6.06	117.37	121.00
3	A	3061	ARG	NE-CZ-NH1	-5.95	117.33	120.30
3	A	3022	ARG	NE-CZ-NH1	-5.94	117.33	120.30
1	C	1399	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	C	1283	ASP	N-CA-C	5.61	126.14	111.00
3	A	3088	ASP	CB-CG-OD1	5.59	123.33	118.30
1	C	1501	VAL	CA-CB-CG2	-5.59	102.51	110.90
2	B	2053	ARG	NE-CZ-NH2	-5.55	117.52	120.30
3	A	3048	ARG	NE-CZ-NH1	-5.54	117.53	120.30
2	B	2031	ASP	CB-CG-OD1	5.52	123.27	118.30
3	A	3022	ARG	NE-CZ-NH2	5.50	123.05	120.30
1	C	1366	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	C	1556	ASP	OD1-CG-OD2	-5.44	112.97	123.30
1	C	1199	VAL	CG1-CB-CG2	5.39	119.52	110.90
1	C	1060	MET	CG-SD-CE	5.37	108.80	100.20
1	C	1306	LEU	CB-CG-CD2	5.33	120.07	111.00
1	C	1133	THR	CA-CB-CG2	-5.33	104.94	112.40
1	C	1319	CYS	N-CA-CB	5.33	120.19	110.60
1	C	1563	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	C	1271	PHE	CB-CG-CD1	5.21	124.45	120.80
1	C	1508	ARG	CD-NE-CZ	5.16	130.82	123.60
1	C	1004	ILE	N-CA-C	-5.12	97.18	111.00
1	C	1136	HIS	CA-CB-CG	-5.12	104.90	113.60
1	C	1339	ARG	CD-NE-CZ	5.10	130.74	123.60
1	C	1318	VAL	CA-CB-CG2	-5.10	103.25	110.90
1	C	1431	ASP	CB-CG-OD1	5.08	122.87	118.30
1	C	1304	TYR	CB-CG-CD2	-5.04	117.98	121.00
1	C	1564	TYR	CZ-CE2-CD2	-5.02	115.28	119.80

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	C	4161	0	4138	50	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	785	0	772	19	0
3	A	776	0	804	14	0
4	C	2	0	0	0	0
5	A	70	0	0	3	0
5	B	76	0	0	1	0
5	C	402	0	0	3	1
All	All	6272	0	5714	78	1

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 7.

All (78) 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:C:1043:VAL:HB	1:C:1051[B]:ILE:HD11	1.48	0.94
1:C:1131:ILE:HD13	1:C:1153:THR:HB	1.51	0.92
2:B:2072:GLU:H	2:B:2075:GLN:HE21	0.93	0.92
1:C:1201:GLN:HE21	1:C:1203:ASP:H	1.13	0.91
2:B:2072:GLU:H	2:B:2075:GLN:NE2	1.70	0.89
2:B:2072:GLU:N	2:B:2075:GLN:HE21	1.78	0.81
3:A:3061:ARG:HD3	3:A:3096:HIS:O	1.84	0.77
2:B:2050:LYS:HE2	5:B:558:HOH:O	1.86	0.75
3:A:3080:ILE:HD12	3:A:3098:PRO:HG3	1.70	0.73
1:C:1043:VAL:CB	1:C:1051[B]:ILE:HD11	2.19	0.72
1:C:1361[B]:SER:OG	1:C:1367:VAL:HB	1.89	0.72
2:B:2024:VAL:CG2	2:B:2055:GLN:HG2	2.20	0.71
3:A:3029:LYS:HE3	3:A:3069:VAL:O	1.90	0.71
3:A:3061:ARG:NH1	5:A:576:HOH:O	2.28	0.67
1:C:1090:ASP:HB3	5:C:516:HOH:O	1.99	0.62
3:A:3061:ARG:HD2	5:A:576:HOH:O	1.99	0.61
1:C:1043:VAL:CG1	1:C:1051[B]:ILE:HD11	2.30	0.61
1:C:1296:SER:HB3	1:C:1356:LEU:HB2	1.83	0.60
1:C:1199:VAL:HG12	1:C:1205:LEU:HD21	1.83	0.60
3:A:3012:LEU:C	3:A:3012:LEU:HD23	2.22	0.60
2:B:2076:LYS:HG2	2:B:2076:LYS:O	2.02	0.59
1:C:1304:TYR:OH	1:C:1339:ARG:HG2	2.04	0.58
1:C:1467:THR:OG1	3:A:3081:GLN:NE2	2.33	0.58
1:C:1131:ILE:HD13	1:C:1153:THR:CB	2.32	0.58
1:C:1395:ASN:HD22	1:C:1397:ASN:H	1.51	0.57
1:C:1022:ARG:O	2:B:2004:GLY:HA2	2.03	0.57
1:C:1390:GLU:HB2	1:C:1398:PHE:CD1	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2024:VAL:HG21	2:B:2055:GLN:HG2	1.87	0.56
2:B:2024:VAL:HG22	2:B:2055:GLN:OE1	2.06	0.55
1:C:1475:PRO:HB2	5:C:411:HOH:O	2.06	0.55
1:C:1233[B]:LEU:HD12	1:C:1243:VAL:HG21	1.89	0.54
3:A:3008:LYS:O	3:A:3011:LEU:HB2	2.08	0.53
1:C:1003:ASN:HA	2:B:2013:ILE:O	2.08	0.52
1:C:1465:ILE:HG13	1:C:1468:PRO:HD3	1.90	0.52
1:C:1390:GLU:N	1:C:1390:GLU:OE1	2.40	0.52
1:C:1246:HIS:CE1	1:C:1278:GLY:HA3	2.46	0.50
3:A:3100:ILE:HD12	5:A:295:HOH:O	2.12	0.50
1:C:1540:THR:OG1	1:C:1542:GLU:HG3	2.12	0.50
3:A:3077:ILE:HG21	3:A:3080:ILE:HD11	1.94	0.49
3:A:3100:ILE:HD12	3:A:3100:ILE:OXT	2.12	0.49
1:C:1080:GLY:HA2	1:C:1401:LYS:HE2	1.95	0.48
1:C:1300:PRO:HG3	1:C:1364:MET:O	2.14	0.48
2:B:2045:VAL:HG21	2:B:2049:LEU:HD12	1.95	0.47
1:C:1505:LEU:O	1:C:1506:ASN:ND2	2.48	0.47
1:C:1020:LYS:HB2	2:B:2007:HIS:HB3	1.98	0.46
3:A:3051:LYS:HD3	3:A:3056:LEU:HD21	1.96	0.46
1:C:1515:LYS:HD3	5:C:264:HOH:O	2.16	0.45
1:C:1154:MET:O	1:C:1191:ILE:HA	2.16	0.45
1:C:1201:GLN:O	1:C:1201:GLN:HG3	2.17	0.45
1:C:1229:ILE:O	1:C:1233[B]:LEU:HD13	2.17	0.45
1:C:1521:GLN:HG3	1:C:1523:ALA:H	1.82	0.45
1:C:1491:PHE:CD2	1:C:1512:ALA:HB3	2.52	0.45
1:C:1068:LEU:HB3	1:C:1088:VAL:HB	1.99	0.45
1:C:1217:KCX:CX	1:C:1219:HIS:HD2	2.30	0.45
1:C:1355:SER:C	1:C:1356:LEU:HD12	2.37	0.44
1:C:1395:ASN:ND2	1:C:1397:ASN:H	2.14	0.44
2:B:2036:VAL:O	2:B:2068:ALA:HB1	2.17	0.44
1:C:1449:LYS:HA	1:C:1449:LYS:HD2	1.82	0.43
1:C:1051[B]:ILE:HD12	1:C:1112:ILE:HD12	2.00	0.43
1:C:1417:ILE:CG2	1:C:1514:VAL:HG12	2.49	0.43
1:C:1186:SER:HB2	1:C:1187[A]:LEU:HD22	1.99	0.42
3:A:3061:ARG:HD2	3:A:3061:ARG:HH11	1.60	0.42
2:B:2059:TYR:HB3	2:B:2081:LEU:HB3	2.01	0.42
2:B:2024:VAL:CG1	2:B:2051:PHE:HB2	2.49	0.42
2:B:2023:ARG:NH1	2:B:2080:GLU:OE2	2.53	0.42
1:C:1277:GLY:HA2	1:C:1364:MET:HE2	2.02	0.42
1:C:1173:GLY:HA2	1:C:1174:PRO:HD3	1.86	0.41
1:C:1359:SER:O	1:C:1365:GLY:HA3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2024:VAL:HG22	2:B:2055:GLN:HG2	2.01	0.41
1:C:1158:GLY:HA3	1:C:1165:THR:HG23	2.03	0.41
1:C:1298:THR:HB	1:C:1360:ASP:HB2	2.03	0.41
1:C:1040:GLY:HA3	2:B:2016:ASN:ND2	2.36	0.40
1:C:1417:ILE:HG22	1:C:1514:VAL:HG12	2.03	0.40
3:A:3011:LEU:HD23	3:A:3011:LEU:HA	1.97	0.40
2:B:2028:ASN:HA	2:B:2049:LEU:HD23	2.04	0.40
1:C:1214:ILE:HB	1:C:1514:VAL:HG11	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:527:HOH:O	5:C:527:HOH:O[15_556]	1.21	0.99

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	C	556/567 (98%)	525 (94%)	29 (5%)	2 (0%)	34 15
2	B	99/101 (98%)	93 (94%)	6 (6%)	0	100 100
3	A	98/100 (98%)	97 (99%)	1 (1%)	0	100 100
All	All	753/768 (98%)	715 (95%)	36 (5%)	2 (0%)	41 21

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1364	MET
1	C	1360	ASP

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	C	438/443 (99%)	426 (97%)	12 (3%)	44	20
2	B	78/78 (100%)	73 (94%)	5 (6%)	17	4
3	A	85/85 (100%)	84 (99%)	1 (1%)	71	54
All	All	601/606 (99%)	583 (97%)	18 (3%)	41	16

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	1043	VAL
1	C	1082	VAL
1	C	1104	ILE
1	C	1246	HIS
1	C	1319	CYS
1	C	1339	ARG
1	C	1391	GLU
1	C	1395	ASN
1	C	1449	LYS
1	C	1506	ASN
1	C	1508	ARG
1	C	1556	ASP
2	B	2009	LYS
2	B	2050	LYS
2	B	2055	GLN
2	B	2076	LYS
2	B	2077	ARG
3	A	3006	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	1003	ASN
1	C	1107	ASN
1	C	1142	GLN

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Mol	Chain	Res	Type
1	C	1201	GLN
1	C	1362	GLN
1	C	1395	ASN
1	C	1469	GLN
2	B	2016	ASN
2	B	2075	GLN
3	A	3081	GLN
3	A	3097	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)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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	$\#\|Z\| > 2$
1	KCX	C	1217	4,1	9,11,12	1.49	1 (11%)	5,12,14	1.91	1 (20%)

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
1	KCX	C	1217	4,1	-	0/9/10/12	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1217	KCX	OQ1-CX	3.46	1.28	1.21

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	1217	KCX	OQ1-CX-NZ	-3.98	118.78	124.96

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	1217	KCX	1	0

5.5 Carbohydrates (i)

There are no monosaccharides 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 (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	C	555/567 (97%)	-0.71	0	100	100	15, 20, 36, 63	0
2	B	101/101 (100%)	-0.27	0	100	100	19, 25, 51, 64	0
3	A	100/100 (100%)	-0.87	0	100	100	15, 20, 38, 57	0
All	All	756/768 (98%)	-0.67	0	100	100	15, 21, 40, 64	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains (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
1	KCX	C	1217	12/13	0.98	0.10	14,18,22,22	0

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
4	NI	C	4774	1/1	1.00	0.06	21,21,21,21	0
4	NI	C	4775	1/1	1.00	0.05	20,20,20,20	0

6.5 Other polymers (i)

There are no such residues in this entry.