



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:14 AM GMT

PDB ID : 2Q01
Title : Crystal structure of glucuronate isomerase from *Caulobacter crescentus*
Authors : Patskovsky, Y.; Bonanno, J.; Sridhar, V.; Sauder, J.M.; Freeman, J.; Powell, A.; Koss, J.; Groshong, C.; Gheyi, T.; Wasserman, S.R.; Raushel, F.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2007-05-18
Resolution : 2.34 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

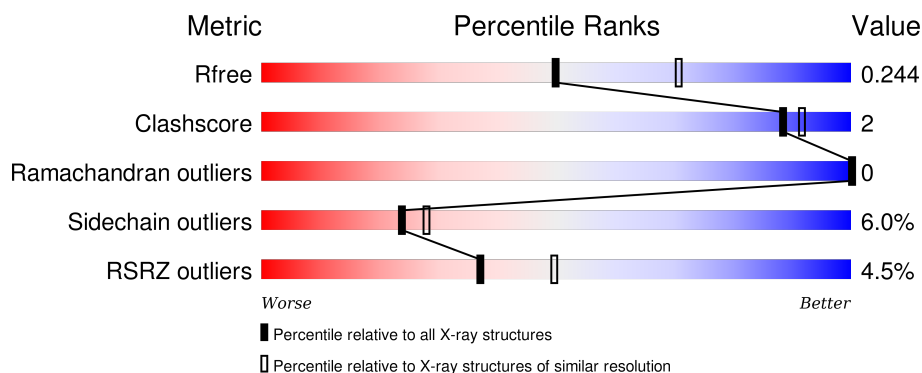
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1406 (2.36-2.32)
Clashscore	102246	1509 (2.36-2.32)
Ramachandran outliers	100387	1490 (2.36-2.32)
Sidechain outliers	100360	1491 (2.36-2.32)
RSRZ outliers	91569	1412 (2.36-2.32)

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	497	<div> <div>5%</div> <div>87%</div> <div>9%</div> <div>• •</div> </div>
1	B	497	<div> <div>4%</div> <div>89%</div> <div>7%</div> <div>• •</div> </div>
1	C	497	<div> <div>3%</div> <div>85%</div> <div>10%</div> <div>• •</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12130 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 Uronate isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	479	Total	C	N	O	S	0	10	0
			3887	2468	685	722	12			
1	B	480	Total	C	N	O	S	0	7	0
			3871	2459	680	721	11			
1	C	480	Total	C	N	O	S	0	7	0
			3881	2465	686	719	11			

There are 33 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	CLONING ARTIFACT	UNP Q9A874
A	0	SER	-	CLONING ARTIFACT	UNP Q9A874
A	1	LEU	-	CLONING ARTIFACT	UNP Q9A874
A	488	GLU	-	CLONING ARTIFACT	UNP Q9A874
A	489	GLY	-	CLONING ARTIFACT	UNP Q9A874
A	490	HIS	-	CLONING ARTIFACT	UNP Q9A874
A	491	HIS	-	CLONING ARTIFACT	UNP Q9A874
A	492	HIS	-	CLONING ARTIFACT	UNP Q9A874
A	493	HIS	-	CLONING ARTIFACT	UNP Q9A874
A	494	HIS	-	CLONING ARTIFACT	UNP Q9A874
A	495	HIS	-	CLONING ARTIFACT	UNP Q9A874
B	-1	MET	-	CLONING ARTIFACT	UNP Q9A874
B	0	SER	-	CLONING ARTIFACT	UNP Q9A874
B	1	LEU	-	CLONING ARTIFACT	UNP Q9A874
B	488	GLU	-	CLONING ARTIFACT	UNP Q9A874
B	489	GLY	-	CLONING ARTIFACT	UNP Q9A874
B	490	HIS	-	CLONING ARTIFACT	UNP Q9A874
B	491	HIS	-	CLONING ARTIFACT	UNP Q9A874
B	492	HIS	-	CLONING ARTIFACT	UNP Q9A874
B	493	HIS	-	CLONING ARTIFACT	UNP Q9A874
B	494	HIS	-	CLONING ARTIFACT	UNP Q9A874
B	495	HIS	-	CLONING ARTIFACT	UNP Q9A874
C	-1	MET	-	CLONING ARTIFACT	UNP Q9A874

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	CLONING ARTIFACT	UNP Q9A874
C	1	LEU	-	CLONING ARTIFACT	UNP Q9A874
C	488	GLU	-	CLONING ARTIFACT	UNP Q9A874
C	489	GLY	-	CLONING ARTIFACT	UNP Q9A874
C	490	HIS	-	CLONING ARTIFACT	UNP Q9A874
C	491	HIS	-	CLONING ARTIFACT	UNP Q9A874
C	492	HIS	-	CLONING ARTIFACT	UNP Q9A874
C	493	HIS	-	CLONING ARTIFACT	UNP Q9A874
C	494	HIS	-	CLONING ARTIFACT	UNP Q9A874
C	495	HIS	-	CLONING ARTIFACT	UNP Q9A874

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

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

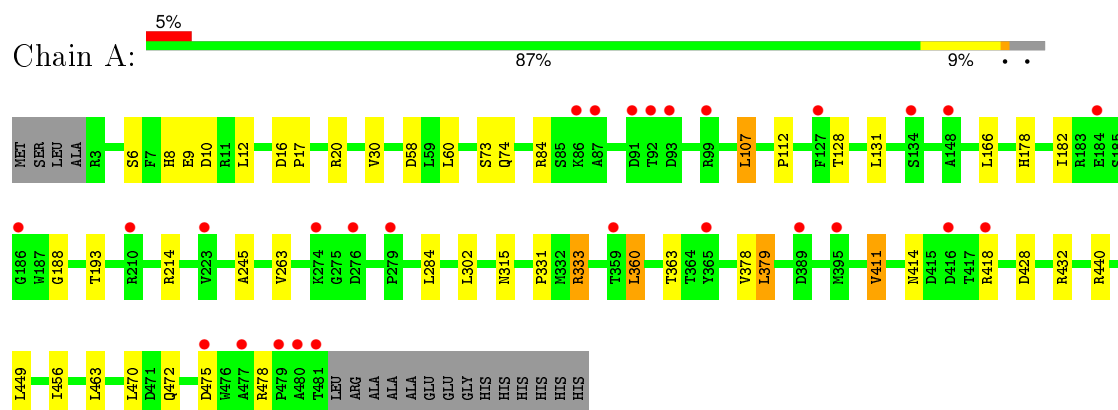
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	151	Total O 151 151	0	0
3	B	188	Total O 188 188	0	0
3	C	149	Total O 149 149	0	0

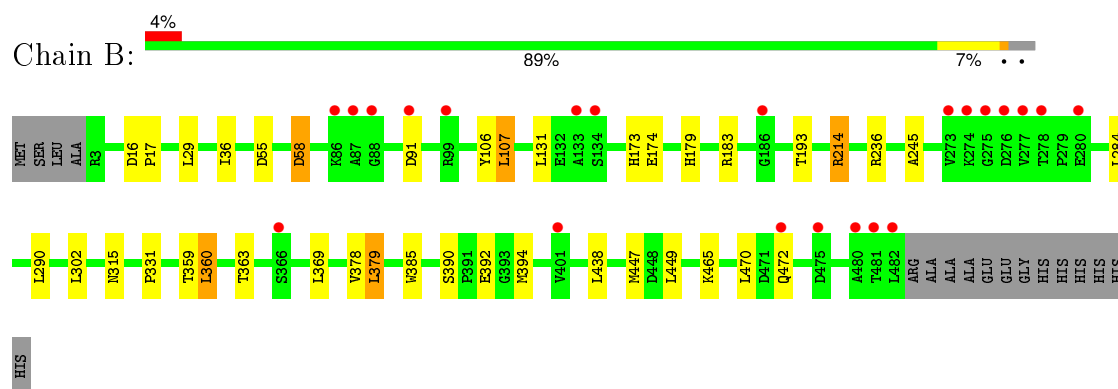
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

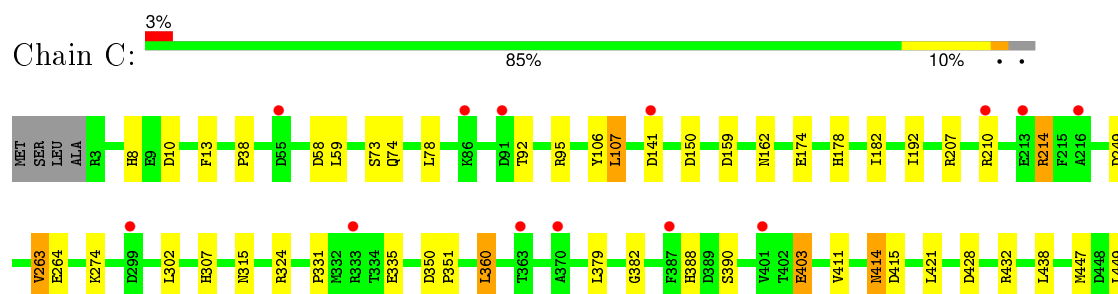
- Molecule 1: Uronate isomerase

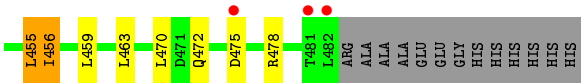


- Molecule 1: Uronate isomerase



- Molecule 1: Uronate isomerase





4 Data and refinement statistics

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	177.33Å 190.18Å 319.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.34 46.44 – 2.34	Depositor EDS
% Data completeness (in resolution range)	91.7 (20.00-2.34) 91.7 (46.44-2.34)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	0.18	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.3.0034	Depositor
R, R_{free}	0.198 , 0.251 0.196 , 0.244	Depositor DCC
R_{free} test set	3035 reflections (3.04%)	DCC
Wilson B-factor (Å ²)	46.0	Xtriage
Anisotropy	0.231	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	2 of 102906 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12130	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.67% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: K

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	0.36	0/4029	0.58	0/5478
1	B	0.36	0/4003	0.56	0/5445
1	C	0.36	0/4013	0.58	0/5457
All	All	0.36	0/12045	0.58	0/16380

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

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	3887	0	3744	20	0
1	B	3871	0	3724	15	0
1	C	3881	0	3740	24	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	151	0	0	0	0
3	B	188	0	0	0	0
3	C	149	0	0	0	0
All	All	12130	0	11208	56	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 2.

All (56) 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:8:HIS:HD2	1:A:10[B]:ASP:H	1.36	0.72
1:A:8:HIS:HD2	1:A:10[A]:ASP:H	1.37	0.72
1:C:8:HIS:HD2	1:C:10[B]:ASP:H	1.48	0.61
1:C:8:HIS:HD2	1:C:10[A]:ASP:H	1.49	0.60
1:A:411:VAL:HG13	1:A:411:VAL:O	2.01	0.60
1:C:331:PRO:HB3	1:C:360:LEU:HB3	1.85	0.59
1:B:179:HIS:O	1:B:183:ARG:HG3	2.02	0.58
1:A:9:GLU:O	1:A:432:ARG:NH2	2.38	0.56
1:B:173:HIS:HB2	1:B:214:ARG:HG3	1.88	0.56
1:A:74:GLN:HG3	1:A:107:LEU:HD21	1.89	0.55
1:A:12:LEU:HD12	1:A:432:ARG:HD3	1.92	0.51
1:C:207:ARG:HE	1:C:324:ARG:HD2	1.76	0.50
1:C:159:ASP:OD2	1:C:478:ARG:NH2	2.45	0.50
1:B:331:PRO:HB3	1:B:360:LEU:HB3	1.95	0.49
1:A:331:PRO:HB3	1:A:360:LEU:HB3	1.95	0.47
1:A:363:THR:HB	1:B:363:THR:HG21	1.96	0.47
1:B:378:VAL:HG23	1:B:379:LEU:HD13	1.96	0.47
1:C:106:TYR:CE2	1:C:107:LEU:CD1	2.97	0.47
1:C:38:PRO:HD2	1:C:414:ASN:HB3	1.97	0.47
1:C:249:ASP:OD1	1:C:307:HIS:ND1	2.45	0.47
1:B:390:SER:O	1:B:394:MET:HG3	2.15	0.47
1:C:428:ASP:O	1:C:432[A]:ARG:HG3	2.16	0.46
1:B:16:ASP:HA	1:B:17:PRO:HD3	1.79	0.46
1:A:178:HIS:O	1:A:182:ILE:HG12	2.16	0.46
1:B:36:ILE:HD13	1:B:465:LYS:HG3	1.98	0.46
1:C:95:ARG:NH1	1:C:141:ASP:OD1	2.48	0.46
1:A:16:ASP:HA	1:A:17:PRO:HD3	1.76	0.45
1:A:74:GLN:HG3	1:A:107:LEU:CD2	2.46	0.45
1:C:59:LEU:HD11	1:C:421:LEU:HD13	1.99	0.45
1:C:263:VAL:CG1	1:C:264:GLU:N	2.79	0.45
1:B:236:ARG:HA	1:B:236:ARG:HD2	1.78	0.45
1:A:378:VAL:HG23	1:A:379:LEU:HD13	1.99	0.44
1:B:360:LEU:HD13	1:B:385:TRP:CE3	2.53	0.44
1:C:106:TYR:CE2	1:C:107:LEU:HD13	2.52	0.43
1:A:188:GLY:HA3	1:A:478:ARG:HG3	2.00	0.43
1:C:382:GLY:HA2	1:C:411:VAL:HG22	2.01	0.43
1:C:447:MET:HE1	1:C:455:LEU:HD23	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:PRO:HD3	1:C:403:GLU:HG3	1.99	0.43
1:A:428:ASP:O	1:A:432:ARG:HG3	2.19	0.42
1:B:290:LEU:HA	1:B:290:LEU:HD23	1.95	0.42
1:C:74:GLN:HG3	1:C:107:LEU:HD21	2.01	0.42
1:C:178:HIS:O	1:C:182:ILE:HG12	2.19	0.42
1:C:350:ASP:HA	1:C:351:PRO:HD3	1.91	0.41
1:B:359:THR:HB	1:B:369:LEU:HD11	2.03	0.41
1:B:55:ASP:HB2	1:B:58:ASP:H	1.86	0.41
1:A:333[A]:ARG:HH11	1:C:335:GLU:HG3	1.86	0.41
1:A:193:THR:HG23	1:A:245:ALA:HB2	2.02	0.41
1:C:210:ARG:O	1:C:214:ARG:HG2	2.21	0.41
1:B:193:THR:HG23	1:B:245:ALA:HB2	2.03	0.41
1:A:10[B]:ASP:OD2	1:A:20:ARG:NH1	2.54	0.41
1:C:456:ILE:HG13	1:C:456:ILE:H	1.61	0.41
1:C:388:HIS:C	1:C:390:SER:H	2.24	0.40
1:C:10[B]:ASP:HB2	1:C:13:PHE:HD2	1.87	0.40
1:B:106:TYR:CE2	1:B:107:LEU:HD13	2.56	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	487/497 (98%)	471 (97%)	16 (3%)	0	100	100
1	B	485/497 (98%)	472 (97%)	13 (3%)	0	100	100
1	C	485/497 (98%)	468 (96%)	17 (4%)	0	100	100
All	All	1457/1491 (98%)	1411 (97%)	46 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	412/415 (99%)	383 (93%)	29 (7%)	19	20
1	B	410/415 (99%)	392 (96%)	18 (4%)	35	44
1	C	410/415 (99%)	382 (93%)	28 (7%)	20	22
All	All	1232/1245 (99%)	1157 (94%)	75 (6%)	24	27

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	VAL
1	A	58	ASP
1	A	60	LEU
1	A	73	SER
1	A	84	ARG
1	A	107	LEU
1	A	128	THR
1	A	131	LEU
1	A	166	LEU
1	A	214	ARG
1	A	263	VAL
1	A	284	LEU
1	A	302	LEU
1	A	315	ASN
1	A	333[A]	ARG
1	A	333[B]	ARG
1	A	360	LEU
1	A	379	LEU
1	A	411	VAL
1	A	414	ASN
1	A	418[A]	ARG
1	A	418[B]	ARG
1	A	440	ARG
1	A	449	LEU
1	A	456	ILE
1	A	463	LEU

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Mol	Chain	Res	Type
1	A	470	LEU
1	A	472	GLN
1	A	475	ASP
1	B	29	LEU
1	B	58	ASP
1	B	91	ASP
1	B	107	LEU
1	B	131	LEU
1	B	174	GLU
1	B	214	ARG
1	B	284	LEU
1	B	302	LEU
1	B	315	ASN
1	B	360	LEU
1	B	379	LEU
1	B	392	GLU
1	B	438	LEU
1	B	447	MET
1	B	449	LEU
1	B	470	LEU
1	B	472	GLN
1	C	58	ASP
1	C	73	SER
1	C	78	LEU
1	C	92	THR
1	C	107	LEU
1	C	150	ASP
1	C	162	ASN
1	C	174	GLU
1	C	192	ILE
1	C	214	ARG
1	C	263	VAL
1	C	274	LYS
1	C	302	LEU
1	C	315	ASN
1	C	360	LEU
1	C	379	LEU
1	C	403	GLU
1	C	414	ASN
1	C	415	ASP
1	C	438	LEU
1	C	449	LEU

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Mol	Chain	Res	Type
1	C	455	LEU
1	C	456	ILE
1	C	459	LEU
1	C	463	LEU
1	C	470	LEU
1	C	472	GLN
1	C	475	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	HIS
1	A	162	ASN
1	A	190	HIS
1	A	270	ASN
1	A	305	GLN
1	A	311	HIS
1	A	315	ASN
1	A	349	ASN
1	A	400	GLN
1	A	414	ASN
1	B	162	ASN
1	B	177	GLN
1	B	270	ASN
1	B	311	HIS
1	B	315	ASN
1	B	349	ASN
1	B	400	GLN
1	B	414	ASN
1	C	8	HIS
1	C	162	ASN
1	C	270	ASN
1	C	305	GLN
1	C	311	HIS
1	C	315	ASN
1	C	349	ASN
1	C	400	GLN
1	C	414	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 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 ⓘ

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	479/497 (96%)	0.33	27 (5%)	28 40	29, 52, 82, 108	0
1	B	480/497 (96%)	0.17	22 (4%)	36 49	28, 46, 74, 109	0
1	C	480/497 (96%)	0.21	16 (3%)	50 61	31, 47, 75, 119	0
All	All	1439/1491 (96%)	0.23	65 (4%)	37 50	28, 48, 77, 119	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	91	ASP	5.8
1	B	482	LEU	4.8
1	A	186	GLY	4.4
1	A	87	ALA	4.4
1	C	91	ASP	4.1
1	A	276[A]	ASP	3.9
1	B	481	THR	3.8
1	A	481	THR	3.7
1	B	276[A]	ASP	3.6
1	A	480	ALA	3.4
1	A	479	PRO	3.2
1	C	481	THR	3.2
1	A	475	ASP	3.2
1	C	482	LEU	3.1
1	B	91	ASP	3.0
1	A	134	SER	2.9
1	B	475	ASP	2.7
1	C	216	ALA	2.6
1	B	277	VAL	2.6
1	C	333[A]	ARG	2.6
1	C	363	THR	2.6
1	A	148	ALA	2.6
1	B	133	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	274	LYS	2.6
1	A	210	ARG	2.5
1	B	278	THR	2.5
1	A	274	LYS	2.5
1	C	86	LYS	2.5
1	C	213	GLU	2.4
1	B	401	VAL	2.4
1	A	99	ARG	2.4
1	C	370	ALA	2.4
1	A	223	VAL	2.4
1	A	359	THR	2.4
1	A	418[A]	ARG	2.4
1	B	186	GLY	2.3
1	A	365	TYR	2.3
1	C	299	ASP	2.3
1	B	472	GLN	2.3
1	B	134	SER	2.3
1	A	93	ASP	2.3
1	B	87	ALA	2.2
1	C	401	VAL	2.2
1	B	366	SER	2.2
1	C	387	PHE	2.2
1	A	279	PRO	2.2
1	B	88	GLY	2.2
1	A	395	MET	2.2
1	A	477	ALA	2.2
1	B	86	LYS	2.2
1	B	99	ARG	2.2
1	A	416	ASP	2.2
1	B	280	GLU	2.2
1	C	141	ASP	2.1
1	A	127	PHE	2.1
1	C	475	ASP	2.1
1	B	480	ALA	2.1
1	C	55	ASP	2.1
1	A	389[A]	ASP	2.1
1	C	210	ARG	2.1
1	A	184	GLU	2.0
1	B	273	VAL	2.0
1	B	275	GLY	2.0
1	A	86	LYS	2.0
1	A	92	THR	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
2	K	B	502	1/1	0.98	0.10	-1.38	37,37,37,37	0
2	K	A	501	1/1	0.99	0.08	-1.96	35,35,35,35	0
2	K	C	503	1/1	0.98	0.09	-1.96	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.