



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 01:15 AM UTC

PDB ID : 4FFX / pdb\_00004ffx  
Title : Structural and Biochemical Characterization of Human Adenylosuccinate Lyase (ADSL) and the R303C ADSL Deficiency Associated Mutation  
Authors : Deaton, M.K.; Ray, S.P.; Capodagli, G.C.; Calkins, L.A.F.; Sawle, L.; Ghosh, K.; Patterson, D.; Pegan, S.D.  
Deposited on : 2012-06-01  
Resolution : 2.70 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](mailto: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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

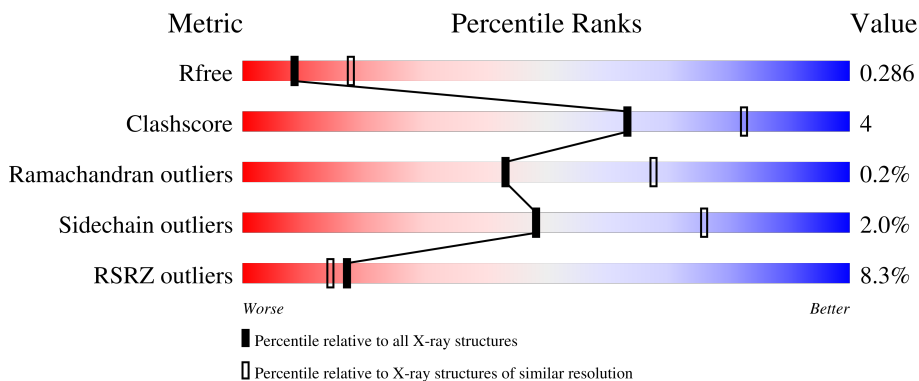
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	487	 10% 84% 9% • 6%
1	B	487	 6% 84% 9% 6%
1	C	487	 9% 80% 11% • 7%
1	D	487	 6% 80% 13% 6%

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 14819 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 Adenylosuccinate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	457	3662	2306	653	678	25	0	2	0
1	B	456	3661	2306	652	678	25	0	2	0
1	C	451	3619	2282	644	667	26	0	3	0
1	D	456	3663	2310	649	679	25	0	4	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P30566
A	-1	SER	-	expression tag	UNP P30566
A	0	HIS	-	expression tag	UNP P30566
A	63	ARG	GLN	conflict	UNP P30566
B	-2	GLY	-	expression tag	UNP P30566
B	-1	SER	-	expression tag	UNP P30566
B	0	HIS	-	expression tag	UNP P30566
B	63	ARG	GLN	conflict	UNP P30566
C	-2	GLY	-	expression tag	UNP P30566
C	-1	SER	-	expression tag	UNP P30566
C	0	HIS	-	expression tag	UNP P30566
C	63	ARG	GLN	conflict	UNP P30566
D	-2	GLY	-	expression tag	UNP P30566
D	-1	SER	-	expression tag	UNP P30566
D	0	HIS	-	expression tag	UNP P30566
D	63	ARG	GLN	conflict	UNP P30566

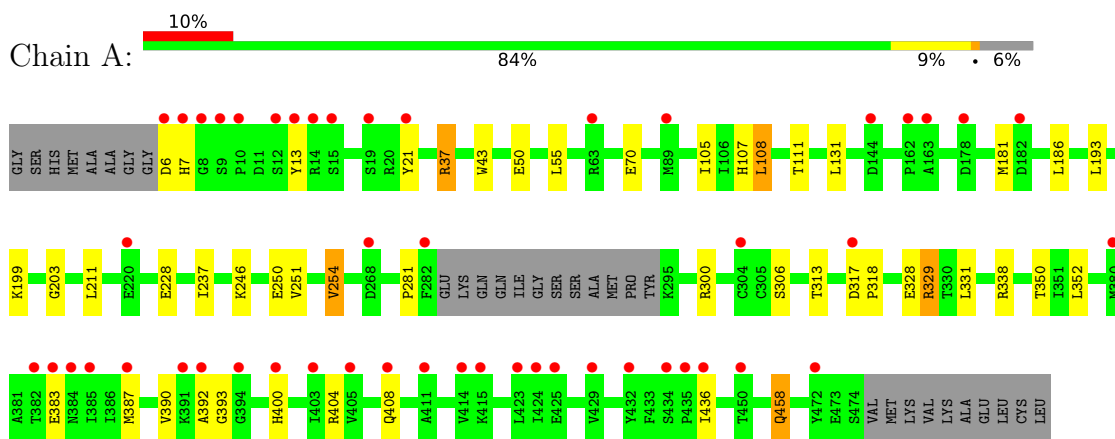
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	67	Total O 67 67	0	0
2	B	45	Total O 45 45	0	0
2	C	38	Total O 38 38	0	0
2	D	64	Total O 64 64	0	0

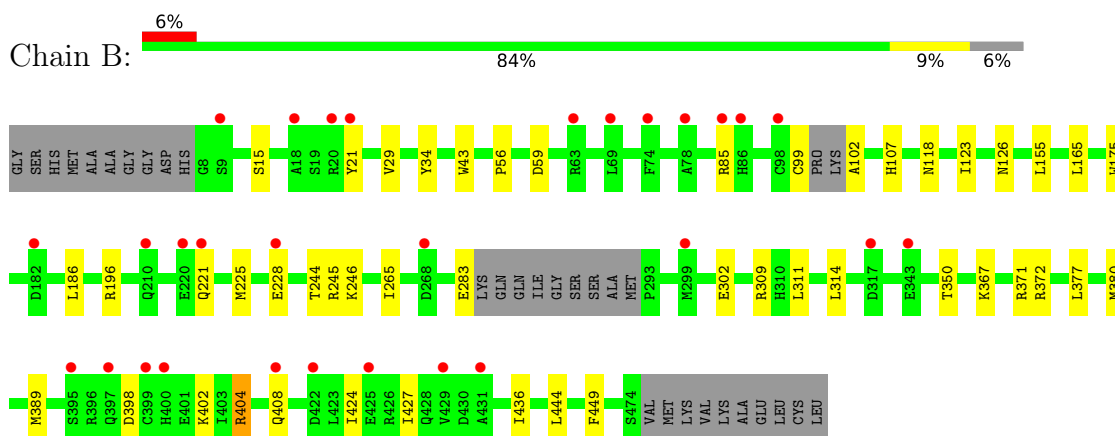
### 3 Residue-property plots [i](#)

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.

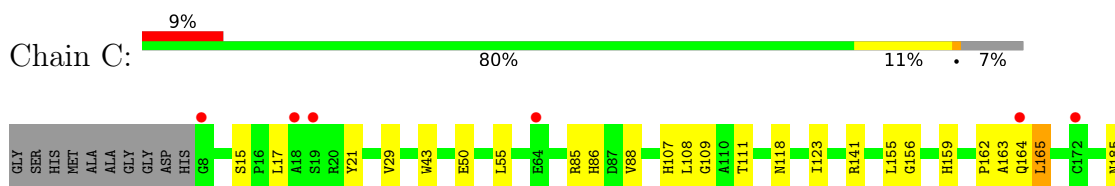
- Molecule 1: Adenylosuccinate lyase

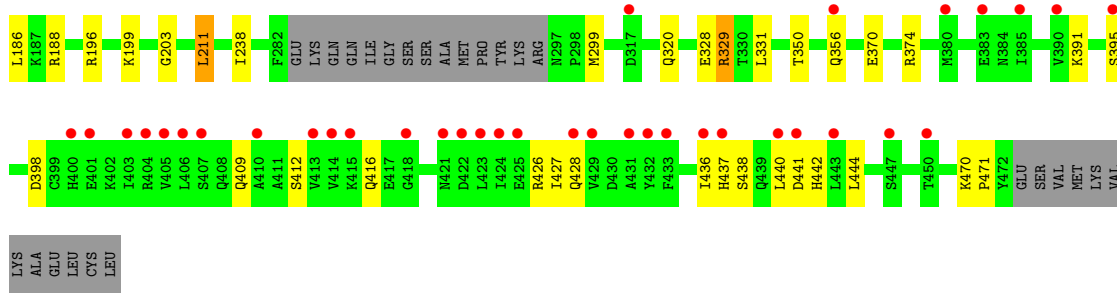


- Molecule 1: Adenylosuccinate lyase

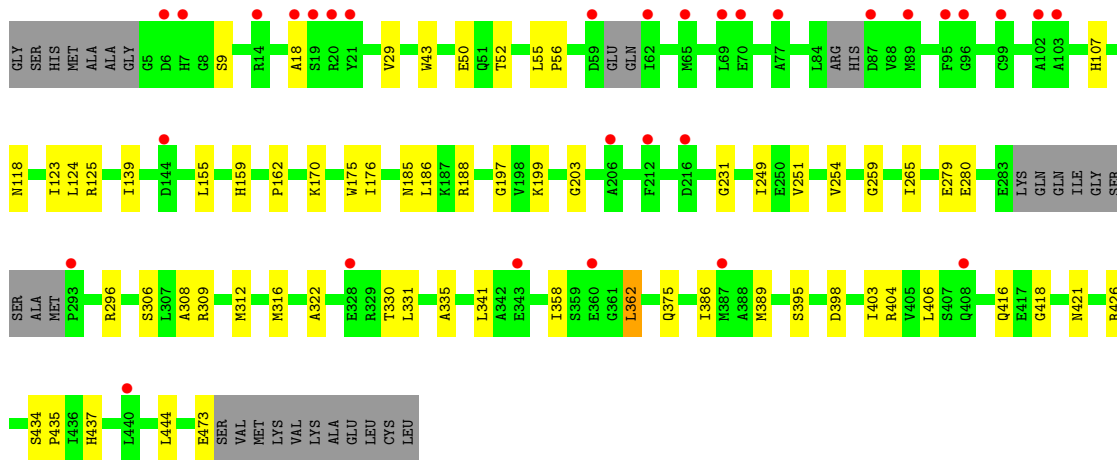
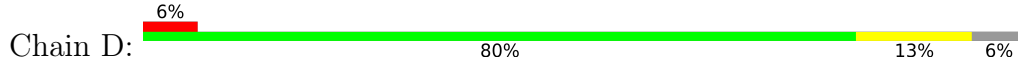


- Molecule 1: Adenylosuccinate lyase





• Molecule 1: Adenylosuccinate lyase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.92Å 105.15Å 214.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 50.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	86.8 (50.00-2.70) 87.2 (50.00-2.70)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.46 (at 2.69Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.243 , 0.292 0.239 , 0.286	Depositor DCC
$R_{free}$ test set	2457 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	25.0	Xtrriage
Anisotropy	0.232	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 24.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	14819	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.38% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	0.43	0/3735	0.82	1/5042 (0.0%)
1	B	0.43	0/3730	0.81	0/5033
1	C	0.42	0/3691	0.81	0/4983
1	D	0.42	0/3738	0.82	2/5043 (0.0%)
All	All	0.43	0/14894	0.81	3/20101 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	237	ILE	N-CA-C	-6.29	106.40	111.81
1	D	280	GLU	CA-C-N	5.28	125.21	119.78
1	D	280	GLU	C-N-CA	5.28	125.21	119.78

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	3662	0	3681	31	0
1	B	3661	0	3676	29	0
1	C	3619	0	3646	38	0
1	D	3663	0	3681	35	0
2	A	67	0	0	1	0
2	B	45	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	38	0	0	1	0
2	D	64	0	0	0	0
All	All	14819	0	14684	116	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 4.

All (116) 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:B:404:ARG:O	1:B:404:ARG:HG3	1.70	0.90
1:C:440:LEU:HB2	1:C:441:ASP:HA	1.56	0.87
1:A:43:TRP:HE1	1:A:107:HIS:HD2	1.27	0.80
1:A:458[B]:GLN:HE21	1:A:458[B]:GLN:HA	1.45	0.78
1:B:15:SER:HB2	1:C:15:SER:HB2	1.67	0.75
1:A:6:ASP:HB2	1:D:9:SER:HB3	1.72	0.72
1:A:43:TRP:HE1	1:A:107:HIS:CD2	2.08	0.72
1:A:328:GLU:O	1:A:329:ARG:HB2	1.93	0.69
1:C:43:TRP:HE1	1:C:107:HIS:CD2	2.12	0.67
1:A:458[B]:GLN:HE21	1:A:458[B]:GLN:CA	2.08	0.66
1:C:43:TRP:HE1	1:C:107:HIS:HD2	1.44	0.65
1:C:165:LEU:H	1:C:165:LEU:HD23	1.61	0.64
1:C:109:GLY:HA3	1:C:211:LEU:HD12	1.80	0.64
1:C:328:GLU:O	1:C:329:ARG:HB2	1.98	0.64
1:B:389:MET:HE3	1:B:436:ILE:HB	1.80	0.63
1:A:181:MET:HE3	1:B:246:LYS:HD3	1.82	0.62
1:D:43:TRP:HE1	1:D:107:HIS:CD2	2.18	0.62
1:B:29:VAL:HG12	1:B:123:ILE:HG23	1.80	0.61
1:C:440:LEU:CB	1:C:441:ASP:HA	2.27	0.59
1:C:395:SER:HB3	1:C:398:ASP:HB2	1.86	0.58
1:A:6:ASP:N	1:D:9:SER:HG	2.02	0.58
1:A:306:SER:HB3	1:D:335:ALA:HB3	1.87	0.57
1:C:111:THR:HG22	2:C:513:HOH:O	2.04	0.57
1:B:85:ARG:NH2	1:C:299[A]:MET:SD	2.78	0.57
1:A:108:LEU:HD12	1:A:211:LEU:HD11	1.87	0.56
1:D:125:ARG:HH12	1:D:231:GLY:C	2.13	0.56
1:C:108:LEU:HG	1:C:211:LEU:HD11	1.89	0.55
1:C:185:ASN:HD22	1:C:188[A]:ARG:HH12	1.53	0.55
1:B:43:TRP:HE1	1:B:107:HIS:CD2	2.24	0.55
1:A:458[B]:GLN:HA	1:A:458[B]:GLN:NE2	2.20	0.55
1:A:50:GLU:HB3	1:A:55:LEU:HD12	1.89	0.54

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:155:LEU:HD21	1:C:162:PRO:HB3	1.91	0.53
1:D:279:GLU:OE1	1:D:296:ARG:HD3	2.07	0.53
1:A:387:MET:O	1:A:390:VAL:HG22	2.09	0.52
1:D:50:GLU:HB3	1:D:55:LEU:HD12	1.90	0.52
1:C:440:LEU:HA	1:C:442:HIS:H	1.74	0.52
1:A:6:ASP:HB2	1:D:9:SER:CB	2.39	0.51
1:A:199:LYS:HB3	1:A:203:GLY:HA2	1.92	0.51
1:A:7:HIS:H	1:A:7:HIS:CD2	2.28	0.51
1:C:441:ASP:HB2	1:C:444:LEU:HB2	1.92	0.51
1:D:118:ASN:HD21	1:D:197:GLY:H	1.59	0.50
1:D:309:ARG:HA	1:D:312:MET:HE3	1.93	0.50
1:C:470:LYS:HB3	1:C:471:PRO:HD3	1.93	0.50
1:A:251:VAL:O	1:A:254:VAL:HG22	2.13	0.49
1:A:392:ALA:N	1:A:393:GLY:HA2	2.27	0.49
1:B:398:ASP:O	1:B:402:LYS:HG2	2.13	0.49
1:B:311:LEU:HA	1:B:314:LEU:HD12	1.95	0.48
1:A:246:LYS:O	1:A:250:GLU:HG2	2.14	0.48
1:B:371:ARG:HD2	1:D:418:GLY:O	2.14	0.48
1:C:141:ARG:NH2	1:C:356:GLN:HG2	2.28	0.48
1:A:111:THR:HG22	2:A:520:HOH:O	2.13	0.47
1:B:21:TYR:HB3	1:B:350:THR:HG21	1.96	0.47
1:D:185:ASN:HD22	1:D:188[A]:ARG:HH12	1.63	0.47
1:A:181:MET:HE1	1:B:244:THR:HG21	1.97	0.47
1:A:21:TYR:HB3	1:A:350:THR:HG21	1.97	0.47
1:B:309:ARG:HD3	1:C:320:GLN:HB2	1.96	0.46
1:C:412:SER:HA	1:C:416:GLN:HG2	1.96	0.46
1:A:313:THR:HG23	1:D:316:MET:HB2	1.98	0.46
1:B:427:ILE:HG22	1:B:427:ILE:O	2.15	0.46
1:C:156:GLY:H	1:C:165:LEU:HA	1.80	0.46
1:C:185:ASN:ND2	1:C:188[A]:ARG:HH12	2.12	0.46
1:D:175:TRP:HB3	1:D:265:ILE:HG12	1.98	0.46
1:A:400:HIS:O	1:A:404:ARG:HB2	2.16	0.46
1:B:404:ARG:O	1:B:404:ARG:CG	2.54	0.46
1:C:199:LYS:HB3	1:C:203:GLY:HA2	1.98	0.46
1:C:370:GLU:HB3	1:C:374:ARG:NH2	2.30	0.46
1:D:358:ILE:O	1:D:362:LEU:HD13	2.16	0.46
1:B:367:LYS:HG3	1:D:416:GLN:O	2.16	0.46
1:C:118:ASN:ND2	1:C:196:ARG:HB3	2.32	0.45
1:C:50:GLU:HB3	1:C:55:LEU:HD12	1.99	0.45
1:D:406:LEU:HD22	1:D:426:ARG:HB3	1.99	0.45
1:B:175:TRP:HB3	1:B:265:ILE:HG12	1.99	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:389:MET:HE1	1:D:435:PRO:HD2	1.98	0.44
1:D:199:LYS:HB3	1:D:203:GLY:HA2	1.99	0.44
1:A:131:LEU:HD22	1:A:352:LEU:HD12	2.00	0.44
1:A:281:PRO:O	1:A:300:ARG:NH2	2.50	0.44
1:C:238:ILE:HD12	1:D:170:LYS:HE3	1.99	0.44
1:B:118:ASN:HD21	1:B:196:ARG:HB3	1.82	0.44
1:B:118:ASN:ND2	1:B:196:ARG:HB3	2.32	0.44
1:B:165:LEU:HD12	1:B:449:PHE:HB2	1.99	0.44
1:C:86:HIS:HE1	1:C:88:VAL:HB	1.83	0.44
1:D:375:GLN:HE22	1:D:421:ASN:H	1.66	0.44
1:B:302:GLU:OE1	1:D:159:HIS:HD2	2.00	0.43
1:C:159:HIS:HD2	1:D:331:LEU:HG	1.82	0.43
1:B:221:GLN:O	1:B:225:MET:HG3	2.18	0.43
1:C:17:LEU:HA	1:C:21:TYR:HB2	1.99	0.43
1:C:118:ASN:HD21	1:C:196:ARG:HB3	1.83	0.43
1:B:99:CYS:HB3	1:B:102:ALA:HB3	1.99	0.43
1:A:37:ARG:HD3	1:A:70:GLU:OE1	2.18	0.43
1:C:29:VAL:HG12	1:C:123:ILE:HG23	2.01	0.43
1:A:21:TYR:HB3	1:A:350:THR:CG2	2.49	0.43
1:B:424:ILE:HD12	1:B:444:LEU:HD11	2.01	0.43
1:D:29:VAL:HG12	1:D:123:ILE:HG23	2.00	0.42
1:B:155:LEU:HD22	1:B:372:ARG:HG2	2.00	0.42
1:B:377:LEU:HA	1:B:380:MET:HE3	1.99	0.42
1:D:249:ILE:HD11	1:D:322:ALA:HB3	2.01	0.42
1:D:434:SER:HA	1:D:437:HIS:ND1	2.35	0.42
1:C:428:GLN:HG3	1:C:437:HIS:HE1	1.84	0.42
1:C:21:TYR:HB3	1:C:350:THR:HG21	2.02	0.42
1:D:124:LEU:HD21	1:D:341:LEU:HD22	2.02	0.42
1:D:43:TRP:HE1	1:D:107:HIS:HD2	1.66	0.41
1:D:259:GLY:O	1:D:308:ALA:HB1	2.20	0.41
1:B:283:GLU:HG3	1:C:85:ARG:HH22	1.86	0.41
1:A:13:TYR:HD2	1:D:18:ALA:HB2	1.86	0.41
1:D:386:ILE:HD11	1:D:403:ILE:HG21	2.03	0.41
1:B:34:TYR:OH	1:B:126:ASN:ND2	2.54	0.41
1:D:139:ILE:HG23	1:D:176:ILE:HG23	2.03	0.41
1:C:409:GLN:HB2	1:C:426:ARG:NH1	2.36	0.41
1:D:251:VAL:O	1:D:254:VAL:HG22	2.21	0.41
1:D:155:LEU:HD21	1:D:162:PRO:HB3	2.03	0.40
1:C:163:ALA:O	1:C:164:GLN:C	2.64	0.40
1:C:437:HIS:HA	1:C:438:SER:HA	1.74	0.40
1:A:317:ASP:CB	1:A:318:PRO:HD3	2.51	0.40

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:ARG:HD3	1:B:245:ARG:HA	1.93	0.40
1:D:395:SER:HB3	1:D:398:ASP:HB2	2.03	0.40
1:A:105:ILE:HA	1:A:108:LEU:HD23	2.02	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	455/487 (93%)	443 (97%)	11 (2%)	1 (0%)	43 68
1	B	452/487 (93%)	433 (96%)	18 (4%)	1 (0%)	43 68
1	C	450/487 (92%)	434 (96%)	15 (3%)	1 (0%)	43 68
1	D	452/487 (93%)	438 (97%)	13 (3%)	1 (0%)	43 68
All	All	1809/1948 (93%)	1748 (97%)	57 (3%)	4 (0%)	43 68

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	329	ARG
1	C	329	ARG
1	B	56	PRO
1	D	56	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	403/423 (95%)	390 (97%)	13 (3%)	34	64
1	B	402/423 (95%)	397 (99%)	5 (1%)	63	84
1	C	398/423 (94%)	391 (98%)	7 (2%)	51	78
1	D	403/423 (95%)	395 (98%)	8 (2%)	48	76
All	All	1606/1692 (95%)	1573 (98%)	33 (2%)	48	75

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ARG
1	A	108	LEU
1	A	186	LEU
1	A	193	LEU
1	A	228	GLU
1	A	254	VAL
1	A	331	LEU
1	A	338	ARG
1	A	383	GLU
1	A	408	GLN
1	A	436	ILE
1	A	458[A]	GLN
1	A	458[B]	GLN
1	B	59	ASP
1	B	186	LEU
1	B	228	GLU
1	B	404	ARG
1	B	408	GLN
1	C	165	LEU
1	C	186	LEU
1	C	211	LEU
1	C	331	LEU
1	C	391	LYS
1	C	427	ILE
1	C	436	ILE
1	D	52	THR
1	D	186	LEU
1	D	306	SER
1	D	330	THR
1	D	362	LEU
1	D	404	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	D	444	LEU
1	D	473	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	HIS
1	A	41	GLN
1	A	68	ASN
1	A	107	HIS
1	A	118	ASN
1	A	126	ASN
1	A	164	GLN
1	A	185	ASN
1	A	217	HIS
1	A	221	GLN
1	A	357	ASN
1	A	408	GLN
1	A	439	GLN
1	B	41	GLN
1	B	86	HIS
1	B	107	HIS
1	B	118	ASN
1	B	126	ASN
1	B	177	GLN
1	B	184	GLN
1	B	185	ASN
1	B	210	GLN
1	B	297	ASN
1	C	68	ASN
1	C	93	HIS
1	C	107	HIS
1	C	118	ASN
1	C	126	ASN
1	C	159	HIS
1	C	184	GLN
1	C	185	ASN
1	C	274	ASN
1	C	325	GLN
1	C	357	ASN
1	C	384	ASN
1	C	437	HIS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	C	458	GLN
1	D	7	HIS
1	D	41	GLN
1	D	107	HIS
1	D	118	ASN
1	D	126	ASN
1	D	159	HIS
1	D	177	GLN
1	D	185	ASN
1	D	210	GLN
1	D	241	GLN
1	D	356	GLN
1	D	375	GLN
1	D	408	GLN
1	D	416	GLN

### 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 oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	457/487 (93%)	0.74	49 (10%) 11 9	6, 23, 64, 70	2 (0%)
1	B	456/487 (93%)	0.64	29 (6%) 25 22	4, 22, 54, 57	2 (0%)
1	C	451/487 (92%)	0.71	42 (9%) 14 12	5, 23, 69, 71	3 (0%)
1	D	456/487 (93%)	0.69	31 (6%) 23 20	6, 24, 57, 64	4 (0%)
All	All	1820/1948 (93%)	0.70	151 (8%) 17 14	4, 23, 62, 71	11 (0%)

All (151) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	431	ALA	4.4
1	B	429	VAL	4.1
1	C	431	ALA	3.9
1	C	405	VAL	3.9
1	D	96	GLY	3.8
1	C	406	LEU	3.7
1	C	425	GLU	3.7
1	C	432	TYR	3.6
1	A	394	GLY	3.6
1	C	429	VAL	3.6
1	B	182	ASP	3.5
1	B	18	ALA	3.5
1	A	19	SER	3.5
1	D	7	HIS	3.4
1	B	395	SER	3.4
1	D	99	CYS	3.4
1	A	21	TYR	3.4
1	D	19[A]	SER	3.3
1	C	390	VAL	3.3
1	C	424	ILE	3.3
1	C	414	VAL	3.2

*Continued on next page...*

*Continued from previous page...*

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	383	GLU	3.2
1	A	434	SER	3.2
1	C	19	SER	3.2
1	D	18	ALA	3.2
1	D	77	ALA	3.2
1	A	7	HIS	3.1
1	C	401	GLU	3.1
1	B	78	ALA	3.1
1	D	6	ASP	3.1
1	D	328	GLU	3.1
1	C	407	SER	3.1
1	D	343	GLU	3.0
1	A	405	VAL	3.0
1	A	162	PRO	3.0
1	C	433	PHE	3.0
1	D	360	GLU	2.9
1	B	9	SER	2.9
1	A	429	VAL	2.9
1	C	428	GLN	2.9
1	A	9	SER	2.9
1	D	95	PHE	2.9
1	D	65	MET	2.8
1	A	415	LYS	2.8
1	C	400	HIS	2.8
1	A	182	ASP	2.8
1	A	382	THR	2.8
1	A	14	ARG	2.8
1	C	423	LEU	2.7
1	B	85	ARG	2.7
1	A	391	LYS	2.7
1	A	268	ASP	2.7
1	B	21	TYR	2.7
1	C	450	THR	2.7
1	B	397	GLN	2.6
1	A	432	TYR	2.6
1	C	418	GLY	2.6
1	D	293	PRO	2.6
1	B	221	GLN	2.6
1	B	317	ASP	2.6
1	A	13	TYR	2.6
1	B	220	GLU	2.6
1	A	435	PRO	2.6

*Continued on next page...*

*Continued from previous page...*

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	380	MET	2.5
1	C	440	LEU	2.5
1	A	220	GLU	2.5
1	C	422	ASP	2.5
1	C	404	ARG	2.5
1	C	443	LEU	2.5
1	D	387	MET	2.5
1	C	415	LYS	2.5
1	D	14	ARG	2.5
1	A	8	GLY	2.5
1	C	8	GLY	2.5
1	A	12	SER	2.5
1	A	15	SER	2.5
1	A	392	ALA	2.5
1	A	178	ASP	2.4
1	A	414	VAL	2.4
1	B	69	LEU	2.4
1	B	425	GLU	2.4
1	B	210	GLN	2.4
1	D	21	TYR	2.4
1	D	20	ARG	2.4
1	A	400	HIS	2.4
1	B	400	HIS	2.4
1	C	395	SER	2.4
1	A	472	TYR	2.4
1	A	385	ILE	2.4
1	B	408	GLN	2.4
1	C	356	GLN	2.4
1	C	413	VAL	2.4
1	A	163	ALA	2.4
1	D	144	ASP	2.3
1	C	436	ILE	2.3
1	A	403	ILE	2.3
1	C	403	ILE	2.3
1	C	64	GLU	2.3
1	B	228	GLU	2.3
1	C	421	ASN	2.3
1	C	447	SER	2.3
1	A	10	PRO	2.2
1	C	164	GLN	2.2
1	D	89	MET	2.2
1	A	304	CYS	2.2

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	C	317	ASP	2.2
1	D	216	ASP	2.2
1	A	424	ILE	2.2
1	A	436	ILE	2.2
1	D	103	ALA	2.2
1	A	383	GLU	2.2
1	A	450	THR	2.2
1	B	399	CYS	2.2
1	C	437	HIS	2.2
1	B	343	GLU	2.2
1	B	20	ARG	2.2
1	C	410	ALA	2.1
1	A	425	GLU	2.1
1	A	380	MET	2.1
1	A	317	ASP	2.1
1	B	422	ASP	2.1
1	A	282	PHE	2.1
1	B	74	PHE	2.1
1	A	89	MET	2.1
1	A	384	ASN	2.1
1	A	423	LEU	2.1
1	B	299	MET	2.1
1	B	86	HIS	2.1
1	A	6	ASP	2.1
1	D	62	ILE	2.1
1	C	18	ALA	2.1
1	D	59	ASP	2.1
1	D	212	PHE	2.1
1	A	411	ALA	2.1
1	D	102	ALA	2.1
1	A	387	MET	2.1
1	D	440	LEU	2.1
1	B	98	CYS	2.1
1	C	172	CYS	2.1
1	C	385	ILE	2.1
1	A	144	ASP	2.0
1	B	63	ARG	2.0
1	B	268	ASP	2.0
1	C	441	ASP	2.0
1	D	70	GLU	2.0
1	D	408	GLN	2.0
1	A	63	ARG	2.0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	D	87	ASP	2.0
1	D	206	ALA	2.0
1	A	408	GLN	2.0
1	D	69	LEU	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 oligosaccharides in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.