



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:16 AM UTC

PDB ID : 4FLC / pdb_00004flc
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 : 2012-06-14
Resolution : 2.60 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 ⓘ 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

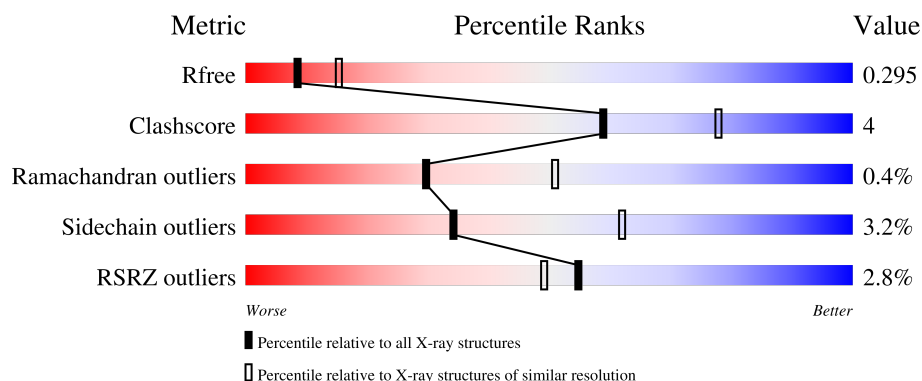
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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))
R_{free}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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 $\leq 5\%$. 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	 2% 78% 15% • 6%
1	B	487	 4% 85% 9% • 5%
1	C	487	 2% 82% 10% 7%
1	D	487	 2% 82% 12% • 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14908 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
1	A	460	Total	C	N	O	S	0	2	0
			3704	2336	658	683	27			
1	B	464	Total	C	N	O	S	0	2	0
			3719	2342	658	692	27			
1	C	453	Total	C	N	O	S	0	3	0
			3632	2290	643	672	27			
1	D	462	Total	C	N	O	S	0	3	0
			3709	2336	660	687	26			

There are 20 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
A	303	CYS	ARG	engineered mutation	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
B	303	CYS	ARG	engineered mutation	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
C	303	CYS	ARG	engineered mutation	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
D	303	CYS	ARG	engineered mutation	UNP P30566

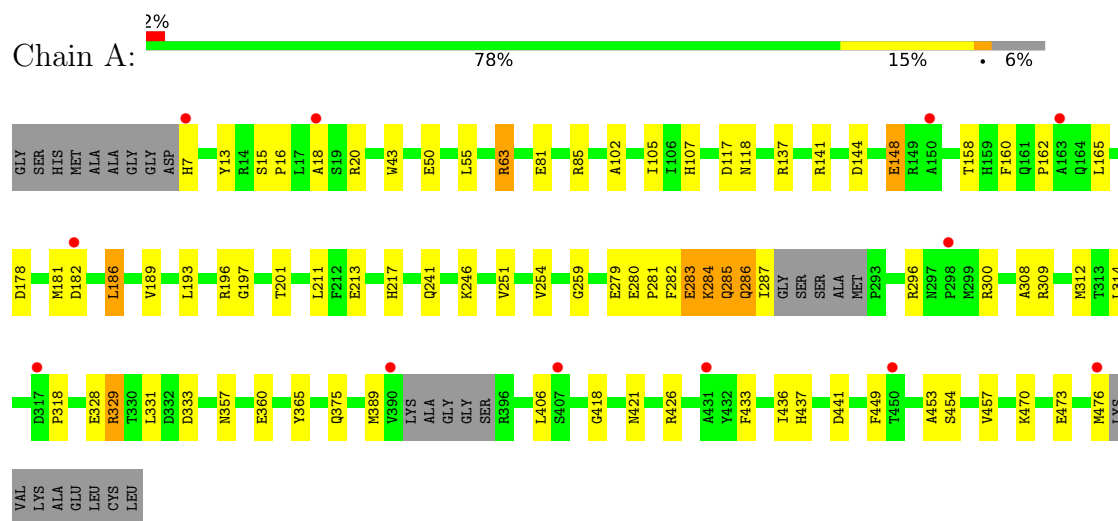
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	36	Total 36	O 36	0	0
2	B	32	Total 32	O 32	0	0
2	C	31	Total 31	O 31	0	0
2	D	45	Total 45	O 45	0	0

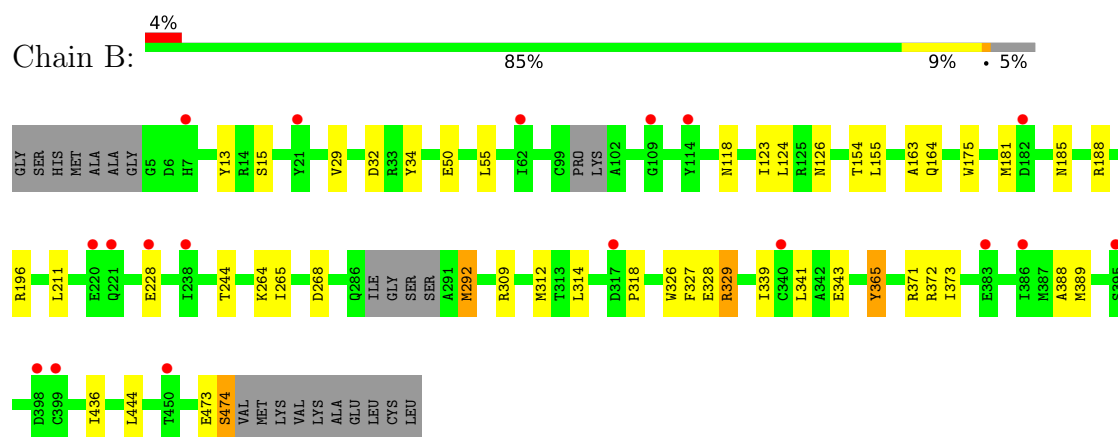
3 Residue-property plots

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.

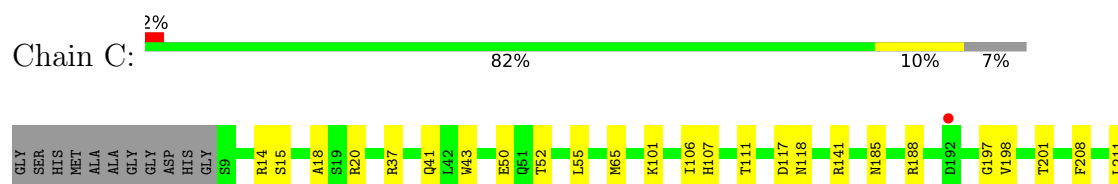
• Molecule 1: Adenylosuccinate lyase

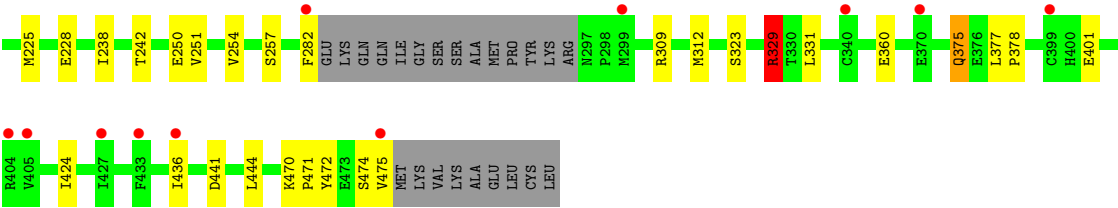


• Molecule 1: Adenylosuccinate lyase

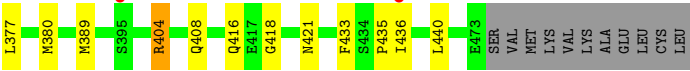
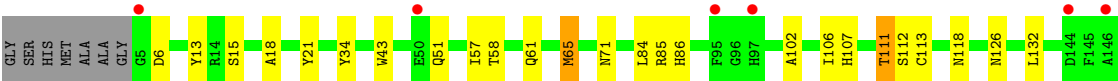
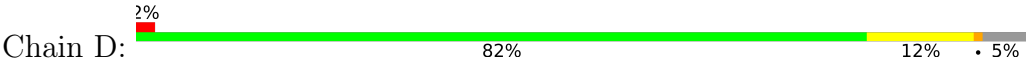


• Molecule 1: Adenylosuccinate lyase





● Molecule 1: Adenylosuccinate lyase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.80Å 105.56Å 217.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.47 – 2.60 47.47 – 2.60	Depositor EDS
% Data completeness (in resolution range)	94.5 (47.47-2.60) 95.1 (47.47-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.29 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.230 , 0.298 0.225 , 0.295	Depositor DCC
R_{free} test set	3136 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14908	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.31% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0/3775	0.85	2/5095 (0.0%)
1	B	0.48	0/3798	0.86	3/5127 (0.1%)
1	C	0.48	0/3704	0.85	0/5002
1	D	0.49	0/3784	0.86	1/5107 (0.0%)
All	All	0.48	0/15061	0.85	6/20331 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	365	TYR	CA-C-N	6.70	126.21	119.24
1	A	365	TYR	C-N-CA	6.70	126.21	119.24
1	D	440	LEU	N-CA-C	5.94	117.83	111.36
1	B	327	PHE	N-CA-C	5.54	116.72	108.31
1	B	365	TYR	CA-C-N	5.35	125.42	119.32
1	B	365	TYR	C-N-CA	5.35	125.42	119.32

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	3704	0	3724	45	0
1	B	3719	0	3733	34	0
1	C	3632	0	3655	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3709	0	3724	37	0
2	A	36	0	0	2	0
2	B	32	0	0	2	0
2	C	31	0	0	2	0
2	D	45	0	0	0	0
All	All	14908	0	14836	131	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 (131) 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:43:TRP:HE1	1:A:107:HIS:HD2	1.10	0.93
1:B:389:MET:HE3	1:B:436:ILE:HB	1.54	0.88
1:C:329:ARG:HG2	1:C:329:ARG:HH11	1.37	0.88
1:A:43:TRP:HE1	1:A:107:HIS:CD2	1.93	0.86
1:A:284:LYS:HG2	1:A:284:LYS:O	1.77	0.85
1:B:473:GLU:O	1:B:474:SER:HB2	1.78	0.84
1:D:375:GLN:HE22	1:D:421:ASN:H	1.27	0.82
1:C:43:TRP:HE1	1:C:107:HIS:CD2	2.00	0.80
1:D:328:GLU:O	1:D:329:ARG:HB2	1.79	0.79
1:A:328:GLU:O	1:A:329:ARG:HB3	1.87	0.73
1:D:404:ARG:O	1:D:408:GLN:HG2	1.88	0.73
1:D:389:MET:HE2	1:D:436:ILE:HG13	1.73	0.69
1:A:281:PRO:O	1:A:300:ARG:NH2	2.25	0.69
1:A:181:MET:HE1	1:B:244[C]:THR:HG21	1.75	0.68
1:C:43:TRP:HE1	1:C:107:HIS:HD2	1.38	0.68
1:A:15:SER:HB2	1:D:15:SER:HB2	1.76	0.68
1:A:181:MET:HE1	1:B:244[A]:THR:HG21	1.75	0.68
1:A:178:ASP:OD1	1:B:244[D]:THR:HG21	1.94	0.67
1:A:118:ASN:HD21	1:A:197:GLY:H	1.42	0.66
1:D:389:MET:CE	1:D:436:ILE:HG13	2.26	0.66
1:A:251:VAL:O	1:A:254:VAL:HG22	1.96	0.66
1:C:185:ASN:HD22	1:C:188[A]:ARG:HH12	1.44	0.64
1:A:178:ASP:HA	1:A:181:MET:HE3	1.81	0.61
1:A:309:ARG:HA	1:A:312:MET:HE3	1.82	0.61
1:B:34:TYR:OH	1:B:126:ASN:ND2	2.34	0.60
1:C:117:ASP:OD2	1:C:242:THR:HB	2.00	0.60
1:D:118:ASN:HD21	1:D:197:GLY:H	1.47	0.60
1:A:13:TYR:HD2	1:D:18:ALA:HB2	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:329:ARG:HG2	1:C:329:ARG:NH1	2.11	0.59
1:D:389:MET:HE2	1:D:436:ILE:CG1	2.32	0.59
1:D:34:TYR:OH	1:D:126:ASN:ND2	2.37	0.58
1:C:37:ARG:HH12	1:C:41:GLN:HE21	1.52	0.56
1:D:21:TYR:HB3	1:D:350:THR:HG21	1.87	0.56
1:A:50:GLU:HB3	1:A:55:LEU:HD12	1.88	0.56
1:D:43:TRP:HE1	1:D:107:HIS:CD2	2.23	0.56
1:B:309:ARG:HA	1:B:312:MET:HE3	1.88	0.56
1:B:388:ALA:HB1	1:B:436:ILE:HD11	1.87	0.56
1:B:124:LEU:HD21	1:B:341:LEU:HD22	1.89	0.55
1:A:283:GLU:O	1:A:284:LYS:HE3	2.07	0.54
1:D:132:LEU:HD21	1:D:186:LEU:HB3	1.89	0.54
1:D:43:TRP:HE1	1:D:107:HIS:HD2	1.54	0.54
1:B:29:VAL:HG12	1:B:123:ILE:HG23	1.90	0.53
1:C:185:ASN:ND2	1:C:188[A]:ARG:HH12	2.07	0.53
1:B:365:TYR:HE1	1:D:416:GLN:HE22	1.56	0.53
1:C:470:LYS:HB3	1:C:471:PRO:HD3	1.90	0.53
1:A:357:ASN:O	1:A:360:GLU:HG3	2.08	0.53
1:B:13:TYR:HD2	1:C:18:ALA:HB2	1.73	0.53
1:B:155:LEU:HD22	1:B:372:ARG:HG2	1.91	0.53
1:A:329:ARG:NH2	1:A:333:ASP:OD1	2.42	0.52
1:B:339:ILE:O	1:B:343:GLU:HG3	2.10	0.52
1:B:50:GLU:O	1:B:55:LEU:HB2	2.10	0.52
1:C:323[B]:SER:OG	1:D:312:MET:SD	2.69	0.51
1:D:84:LEU:O	1:D:85:ARG:HB2	2.09	0.51
1:B:371:ARG:HD2	1:D:418:GLY:O	2.10	0.51
1:C:118:ASN:HD21	1:C:197:GLY:H	1.59	0.50
1:C:111:THR:HG22	2:C:501:HOH:O	2.10	0.50
1:C:238:ILE:HD12	1:D:170:LYS:HE3	1.94	0.50
1:A:279:GLU:HB3	1:A:296:ARG:HH21	1.76	0.49
1:D:65:MET:HE3	1:D:106:ILE:HG12	1.95	0.49
1:A:375:GLN:HE22	1:A:421:ASN:H	1.60	0.49
1:A:217:HIS:HD2	2:A:531:HOH:O	1.96	0.49
1:A:144:ASP:O	1:A:148:GLU:HG3	2.13	0.49
1:A:470:LYS:O	1:A:473:GLU:HG2	2.13	0.49
1:D:317:ASP:CB	1:D:318:PRO:HD3	2.43	0.49
1:C:52:THR:HG21	1:C:225:MET:HE1	1.95	0.48
1:C:378:PRO:HG3	1:C:424:ILE:HD11	1.95	0.48
1:D:51:GLN:HB2	1:D:57:ILE:HB	1.95	0.48
1:C:251:VAL:O	1:C:254:VAL:HG22	2.14	0.48
1:C:375:GLN:HA	1:C:375:GLN:HE21	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:472:TYR:C	1:C:474:SER:H	2.21	0.48
1:D:389:MET:HE1	1:D:435:PRO:HD2	1.94	0.48
1:D:389:MET:CE	1:D:433:PHE:HA	2.44	0.47
1:D:171:ARG:O	1:D:174:LEU:HB2	2.14	0.47
1:A:102:ALA:O	1:A:105:ILE:HG13	2.15	0.47
1:B:175:TRP:HB3	1:B:265:ILE:HG12	1.97	0.47
1:C:309:ARG:HA	1:C:312:MET:HE3	1.96	0.47
1:A:285:GLN:O	1:A:286:GLN:HB3	2.15	0.47
1:C:198:VAL:HB	1:C:208:PHE:CE2	2.50	0.47
1:A:454:SER:HB3	2:A:521:HOH:O	2.15	0.46
1:A:241:GLN:OE1	1:A:329:ARG:HD3	2.15	0.46
1:C:257:SER:HB3	2:C:505:HOH:O	2.16	0.46
1:B:473:GLU:O	1:B:474:SER:CB	2.59	0.45
1:A:389:MET:HE1	1:A:433:PHE:HA	1.99	0.45
1:A:158:THR:HA	1:B:329:ARG:O	2.17	0.45
1:A:16:PRO:O	1:A:20:ARG:HB3	2.17	0.45
1:D:375:GLN:HE22	1:D:421:ASN:N	2.05	0.45
1:A:406:LEU:HD22	1:A:426:ARG:HB3	1.99	0.45
1:A:160:PHE:HB2	1:B:326:TRP:HE3	1.83	0.44
1:C:474:SER:HA	1:C:475:VAL:HA	1.78	0.44
1:C:424:ILE:HD12	1:C:444:LEU:HD13	2.00	0.44
1:D:155:LEU:HD22	1:D:372:ARG:HG2	1.99	0.44
1:A:18:ALA:HB2	1:D:13:TYR:HD2	1.82	0.44
1:B:292:MET:H	1:B:292:MET:HG2	1.62	0.44
1:A:182:ASP:O	1:A:186:LEU:HD12	2.18	0.44
1:B:118:ASN:ND2	1:B:196:ARG:HB3	2.33	0.44
1:C:65:MET:HE3	1:C:106:ILE:HG12	1.99	0.43
1:A:137:ARG:HG2	1:A:141:ARG:HE	1.83	0.43
1:D:309:ARG:HA	1:D:312:MET:HE3	2.01	0.43
1:B:32:ASP:OD2	1:C:14:ARG:NH2	2.52	0.43
1:A:280:GLU:HB3	1:A:300:ARG:NH2	2.34	0.43
1:A:453:ALA:O	1:A:457:VAL:HG23	2.18	0.43
1:B:264:LYS:NZ	1:B:268:ASP:OD2	2.52	0.43
1:B:314:LEU:O	1:B:318:PRO:HD2	2.18	0.43
1:A:165:LEU:HD12	1:A:449:PHE:HB2	2.00	0.43
1:B:15:SER:HB2	1:C:15:SER:HB2	2.00	0.42
1:C:50:GLU:HB3	1:C:55:LEU:HD12	2.01	0.42
1:B:163:ALA:O	1:B:164:GLN:C	2.61	0.42
1:B:188:ARG:NH1	2:B:525:HOH:O	2.52	0.42
1:B:328:GLU:O	1:B:329:ARG:HB2	2.19	0.42
1:D:375:GLN:NE2	1:D:421:ASN:H	2.05	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:377:LEU:HA	1:D:380:MET:HE3	2.01	0.42
1:A:389:MET:HE3	1:A:436:ILE:HB	2.02	0.42
1:D:198:VAL:HB	1:D:208:PHE:CE2	2.55	0.42
1:B:185:ASN:HD22	1:B:188:ARG:HH21	1.66	0.42
1:C:377:LEU:N	1:C:378:PRO:CD	2.83	0.42
1:D:262:VAL:HG11	1:D:351:ILE:HG23	2.02	0.41
1:A:314:LEU:O	1:A:318:PRO:HD2	2.20	0.41
1:A:246:LYS:HD3	1:B:181:MET:HE2	2.02	0.41
1:D:61:GLN:HB3	1:D:102:ALA:HB2	2.02	0.41
1:D:111:THR:CG2	1:D:112:SER:N	2.84	0.41
1:B:13:TYR:CD2	1:C:18:ALA:HB2	2.55	0.41
1:A:63:ARG:HA	1:A:63:ARG:HD3	1.75	0.41
1:D:111:THR:HG22	1:D:113:CYS:H	1.86	0.41
1:D:317:ASP:OD2	1:D:340:CYS:HB3	2.21	0.41
1:C:141:ARG:NE	1:C:360:GLU:HG3	2.36	0.41
1:B:154:THR:HA	1:B:373:ILE:HD11	2.03	0.40
1:B:474:SER:HA	2:B:529:HOH:O	2.20	0.40
1:A:81:GLU:O	1:A:85:ARG:HA	2.22	0.40
1:A:259:GLY:O	1:A:308:ALA:HB1	2.22	0.40
1:A:118:ASN:ND2	1:A:196:ARG:HB3	2.36	0.40
1:C:329:ARG:NH1	1:C:329:ARG:CG	2.83	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	456/487 (94%)	437 (96%)	14 (3%)	5 (1%)	11	25
1	B	462/487 (95%)	445 (96%)	17 (4%)	0	100	100
1	C	452/487 (93%)	436 (96%)	15 (3%)	1 (0%)	43	66
1	D	461/487 (95%)	448 (97%)	11 (2%)	2 (0%)	30	51

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1831/1948 (94%)	1766 (96%)	57 (3%)	8 (0%)	30	51

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	329	ARG
1	D	329	ARG
1	A	213	GLU
1	A	286	GLN
1	C	329	ARG
1	D	332	ASP
1	A	162	PRO
1	A	418	GLY

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	409/423 (97%)	391 (96%)	18 (4%)	25	50
1	B	410/423 (97%)	404 (98%)	6 (2%)	57	80
1	C	401/423 (95%)	388 (97%)	13 (3%)	34	62
1	D	408/423 (96%)	394 (97%)	14 (3%)	32	60
All	All	1628/1692 (96%)	1577 (97%)	51 (3%)	34	63

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	HIS
1	A	63	ARG
1	A	117	ASP
1	A	148	GLU
1	A	186	LEU
1	A	189	VAL
1	A	193	LEU

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Mol	Chain	Res	Type
1	A	201	THR
1	A	211	LEU
1	A	282	PHE
1	A	283	GLU
1	A	284	LYS
1	A	285	GLN
1	A	287	ILE
1	A	331	LEU
1	A	437	HIS
1	A	441	ASP
1	A	476	MET
1	B	211	LEU
1	B	228	GLU
1	B	292	MET
1	B	329	ARG
1	B	444	LEU
1	B	474	SER
1	C	20	ARG
1	C	101	LYS
1	C	201	THR
1	C	211	LEU
1	C	228	GLU
1	C	250	GLU
1	C	282	PHE
1	C	329	ARG
1	C	331	LEU
1	C	375	GLN
1	C	401	GLU
1	C	436	ILE
1	C	441	ASP
1	D	6	ASP
1	D	58	THR
1	D	65	MET
1	D	71	ASN
1	D	86	HIS
1	D	111	THR
1	D	186	LEU
1	D	207	SER
1	D	211	LEU
1	D	220	GLU
1	D	221	GLN
1	D	329	ARG

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Mol	Chain	Res	Type
1	D	334	SER
1	D	404	ARG

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	41	GLN
1	A	97	HIS
1	A	107	HIS
1	A	118	ASN
1	A	126	ASN
1	A	159	HIS
1	A	177	GLN
1	A	184	GLN
1	A	185	ASN
1	A	210	GLN
1	A	221	GLN
1	A	274	ASN
1	A	285	GLN
1	A	336	ASN
1	A	408	GLN
1	A	421	ASN
1	B	7	HIS
1	B	91	HIS
1	B	118	ASN
1	B	126	ASN
1	B	159	HIS
1	B	184	GLN
1	B	185	ASN
1	B	210	GLN
1	B	325	GLN
1	B	400	HIS
1	C	41	GLN
1	C	107	HIS
1	C	118	ASN
1	C	185	ASN
1	C	210	GLN
1	C	356	GLN
1	C	357	ASN
1	C	375	GLN
1	C	458	GLN
1	D	7	HIS

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Mol	Chain	Res	Type
1	D	41	GLN
1	D	71	ASN
1	D	97	HIS
1	D	107	HIS
1	D	118	ASN
1	D	126	ASN
1	D	159	HIS
1	D	184	GLN
1	D	185	ASN
1	D	274	ASN
1	D	356	GLN
1	D	375	GLN
1	D	416	GLN
1	D	439	GLN
1	D	458	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	460/487 (94%)	0.32	12 (2%) 57 51	18, 51, 93, 102	2 (0%)
1	B	464/487 (95%)	0.31	18 (3%) 43 38	17, 53, 83, 84	2 (0%)
1	C	453/487 (93%)	0.32	12 (2%) 57 51	22, 51, 84, 93	3 (0%)
1	D	462/487 (94%)	0.28	10 (2%) 62 57	23, 50, 80, 86	3 (0%)
All	All	1839/1948 (94%)	0.31	52 (2%) 55 49	17, 51, 84, 102	10 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	427	ILE	5.5
1	D	395	SER	3.7
1	A	390	VAL	3.7
1	D	223	ASP	3.4
1	B	398	ASP	3.4
1	D	360	GLU	3.2
1	A	182	ASP	3.2
1	D	5	GLY	3.1
1	B	182	ASP	2.9
1	B	21	TYR	2.9
1	D	144	ASP	2.9
1	C	475	VAL	2.9
1	C	340	CYS	2.8
1	B	317	ASP	2.8
1	C	370	GLU	2.8
1	B	238	ILE	2.8
1	C	299[A]	MET	2.7
1	A	18	ALA	2.7
1	C	433	PHE	2.7
1	B	383	GLU	2.6
1	C	282	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	450	THR	2.6
1	B	386	ILE	2.6
1	B	340	CYS	2.6
1	A	298	PRO	2.5
1	D	95	PHE	2.5
1	C	405	VAL	2.5
1	C	399	CYS	2.4
1	A	407	SER	2.3
1	B	228	GLU	2.3
1	B	450	THR	2.3
1	B	220	GLU	2.3
1	A	163	ALA	2.2
1	B	395	SER	2.2
1	B	7	HIS	2.2
1	D	97	HIS	2.2
1	A	317	ASP	2.2
1	B	62	ILE	2.2
1	D	146	ALA	2.2
1	C	192	ASP	2.2
1	A	150	ALA	2.1
1	A	431	ALA	2.1
1	A	476	MET	2.1
1	B	221	GLN	2.1
1	D	50	GLU	2.1
1	B	109	GLY	2.1
1	B	399	CYS	2.1
1	B	114	TYR	2.1
1	C	436	ILE	2.1
1	D	436	ILE	2.1
1	A	7	HIS	2.0
1	C	404	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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.