



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 09:37 PM UTC

PDB ID : 8YM7 / pdb_00008ym7
Title : Crystal structure of Lysine Specific Demethylase 1 (LSD1) with JH-45
Authors : Zhiyan, D.; Danyan, C.; Hong, J.; Tongchao, L.; Bing, X.
Deposited on : 2024-03-08
Resolution : 2.83 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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

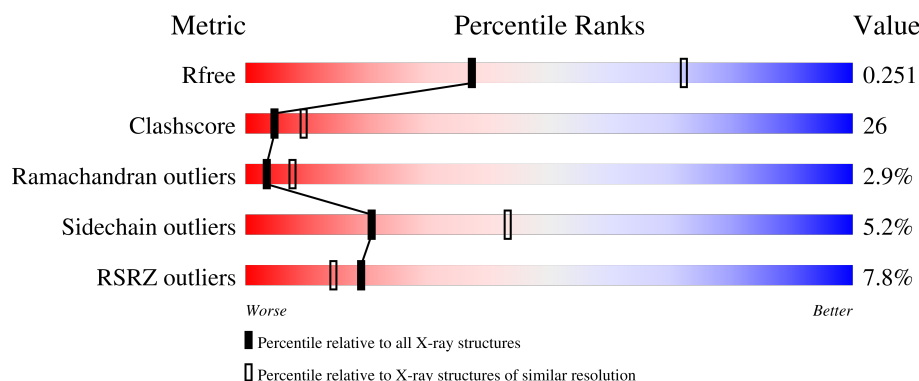
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

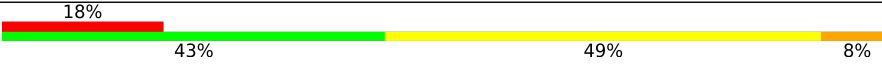
The reported resolution of this entry is 2.83 Å.

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	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	660	
2	B	134	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6331 atoms, of which 23 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 Lysine-specific histone demethylase 1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	660	Total	C	N	O	S	0	0	0
			5162	3288	899	956	19			

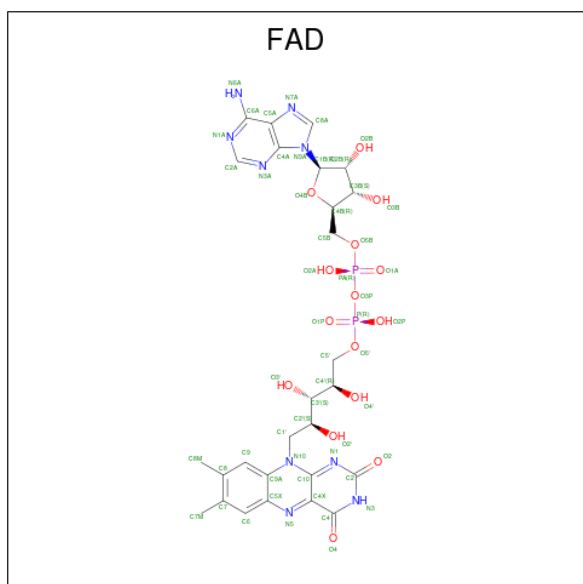
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	173	ALA	GLY	conflict	UNP O60341

- Molecule 2 is a protein called REST corepressor 1.

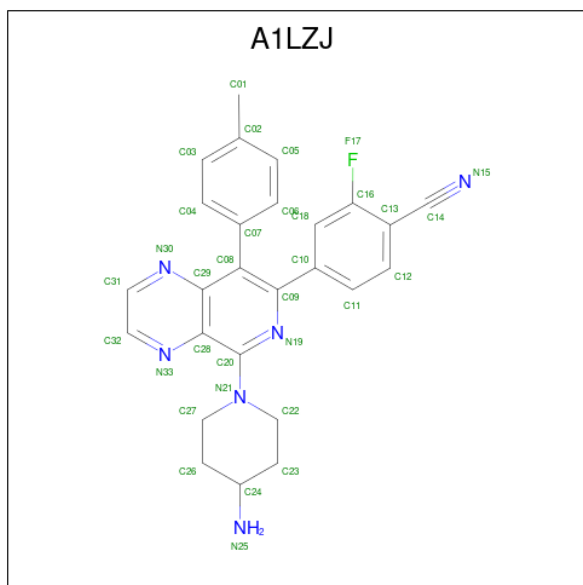
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	134	Total	C	N	O	S	0	0	0
			1060	666	190	201	3			

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 4 is 4-[5-(4-azanylpiperidin-1-yl)-8-(4-methylphenyl)pyrido[3,4-b]pyrazin-7-yl]-2-fluoranyl-benzenecarbonitrile (CCD ID: A1LZJ) (formula: C₂₆H₂₃FN₆) (labeled as "Ligand of Interest" by depositor).

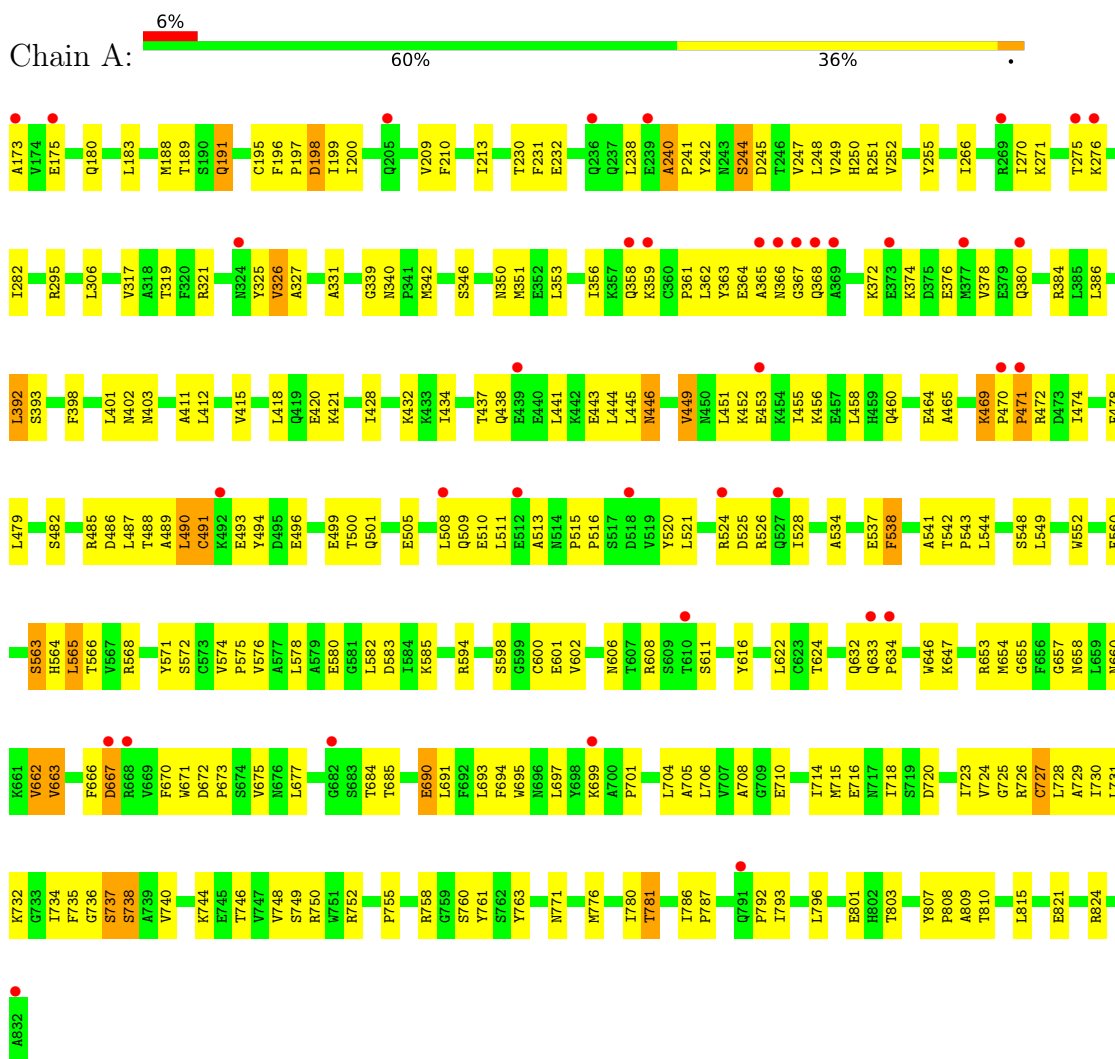


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	F	H	N	0	1
			56	26	1	23	6		

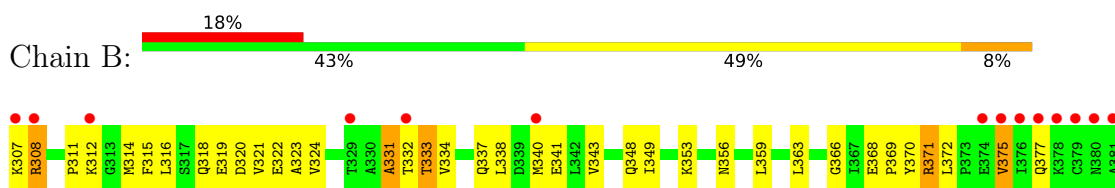
3 Residue-property plots

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: Lysine-specific histone demethylase 1A



• Molecule 2: REST corepressor 1





4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	120.67Å 178.04Å 234.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	117.17 – 2.83 117.17 – 2.83	Depositor EDS
% Data completeness (in resolution range)	99.5 (117.17-2.83) 89.0 (117.17-2.83)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.82 (at 2.82Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.212 , 0.248 0.213 , 0.251	Depositor DCC
R_{free} test set	2000 reflections (3.30%)	wwPDB-VP
Wilson B-factor (Å ²)	57.5	Xtriage
Anisotropy	0.371	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6331	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FAD, A1LZJ

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	0/5274	0.67	3/7157 (0.0%)
2	B	0.36	0/1075	0.59	0/1453
All	All	0.45	0/6349	0.66	3/8610 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	653	ARG	CA-C-N	-5.97	110.06	122.07
1	A	653	ARG	C-N-CA	-5.97	110.06	122.07
1	A	542	THR	N-CA-C	5.35	112.94	108.13

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	469	LYS	Peptide

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	5162	0	5190	260	0
2	B	1060	0	1052	85	0
3	A	53	0	31	5	0
4	A	33	23	0	0	0
All	All	6308	23	6273	321	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 26.

All (321) 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:188:MET:HE3	1:A:200:ILE:HD12	1.34	1.07
1:A:437:THR:HG22	1:A:508:LEU:HD12	1.37	1.06
1:A:566:THR:HG21	1:A:697:LEU:HD22	1.36	1.04
1:A:654:MET:HE1	1:A:776:MET:HG2	1.45	0.98
2:B:311:PRO:HG2	2:B:314:MET:HG3	1.40	0.98
1:A:240:ALA:HB1	1:A:241:PRO:HD2	1.43	0.98
1:A:732:LYS:HG2	1:A:737:SER:HA	1.44	0.98
1:A:366:ASN:HB3	1:A:368:GLN:NE2	1.88	0.89
1:A:240:ALA:HB1	1:A:241:PRO:CD	2.02	0.88
1:A:438:GLN:HE22	2:B:353:LYS:HG3	1.38	0.88
1:A:366:ASN:HB3	1:A:368:GLN:HE22	1.39	0.86
1:A:188:MET:HE3	1:A:200:ILE:CD1	2.05	0.85
1:A:361:PRO:HB2	1:A:363:TYR:HE1	1.41	0.84
1:A:693:LEU:HD12	1:A:694:PHE:H	1.42	0.83
1:A:726:ARG:O	1:A:730:ILE:HG13	1.79	0.82
2:B:411:ASN:O	2:B:412:LYS:HD2	1.79	0.82
1:A:654:MET:CE	1:A:776:MET:HG2	2.09	0.81
2:B:421:PHE:HE1	2:B:434:LEU:HD11	1.45	0.81
1:A:441:LEU:HD23	2:B:356:ASN:HD22	1.45	0.80
1:A:346:SER:CB	1:A:351:MET:HE3	2.11	0.79
1:A:231:PHE:HE1	1:A:249:VAL:HG12	1.48	0.79
1:A:240:ALA:CB	1:A:241:PRO:HD2	2.13	0.79
1:A:446:ASN:ND2	2:B:359:LEU:HD21	1.97	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:438:GLN:NE2	2:B:353:LYS:HG3	1.97	0.78
1:A:356:ILE:O	1:A:358:GLN:HG2	1.84	0.77
1:A:363:TYR:HD2	1:A:734:ILE:HG13	1.48	0.77
1:A:331:ALA:HA	3:A:901:FAD:N5	1.99	0.77
2:B:415:VAL:HG13	2:B:416:GLN:H	1.50	0.76
2:B:307:LYS:HD2	2:B:308:ARG:HE	1.50	0.76
1:A:750:ARG:HG3	1:A:750:ARG:HH11	1.51	0.76
1:A:438:GLN:HG2	1:A:508:LEU:HD11	1.69	0.74
1:A:671:TRP:O	1:A:673:PRO:HD3	1.87	0.74
1:A:331:ALA:HA	3:A:901:FAD:C4X	2.18	0.74
1:A:537:GLU:HG2	1:A:544:LEU:CD1	2.17	0.74
1:A:537:GLU:HG2	1:A:544:LEU:HD13	1.68	0.73
2:B:415:VAL:HG13	2:B:416:GLN:OE1	1.88	0.72
1:A:667:ASP:OD1	1:A:667:ASP:N	2.20	0.72
2:B:322:GLU:OE1	2:B:322:GLU:N	2.17	0.72
1:A:786:ILE:H	1:A:786:ILE:HD12	1.55	0.72
1:A:801:GLU:HG2	1:A:809:ALA:H	1.54	0.71
1:A:734:ILE:HG22	1:A:735:PHE:CD1	2.26	0.71
1:A:456:LYS:HG2	2:B:370:TYR:HE2	1.55	0.71
1:A:666:PHE:O	1:A:701:PRO:HG2	1.90	0.71
1:A:456:LYS:HG2	2:B:370:TYR:CE2	2.26	0.70
1:A:346:SER:HB3	1:A:351:MET:HE3	1.73	0.70
1:A:606:ASN:HD21	1:A:608:ARG:HH21	1.37	0.70
2:B:440:GLU:OE1	2:B:440:GLU:HA	1.90	0.70
1:A:470:PRO:N	1:A:471:PRO:HD2	2.07	0.69
2:B:416:GLN:HA	2:B:419:ASN:HB2	1.75	0.69
1:A:340:ASN:HB2	1:A:560:PHE:CD2	2.27	0.69
1:A:693:LEU:HD12	1:A:694:PHE:N	2.09	0.68
1:A:727:CYS:O	1:A:731:LEU:HD12	1.93	0.68
1:A:441:LEU:HD23	2:B:356:ASN:ND2	2.09	0.68
1:A:446:ASN:HD22	2:B:359:LEU:HD21	1.58	0.67
1:A:465:ALA:HB1	1:A:479:LEU:HD23	1.77	0.67
1:A:750:ARG:HG3	1:A:750:ARG:NH1	2.08	0.67
1:A:510:GLU:HG2	1:A:511:LEU:HD23	1.78	0.66
1:A:690:GLU:OE2	1:A:726:ARG:NH1	2.28	0.66
1:A:245:ASP:OD1	1:A:247:VAL:HG12	1.94	0.66
1:A:740:VAL:HG12	1:A:740:VAL:O	1.96	0.66
1:A:485:ARG:HG3	1:A:485:ARG:HH11	1.59	0.66
1:A:241:PRO:O	1:A:244:SER:HB3	1.95	0.66
1:A:458:LEU:HB3	1:A:487:LEU:HD13	1.77	0.65
2:B:337:GLN:HA	2:B:340:MET:HE2	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:ALA:CB	1:A:241:PRO:CD	2.72	0.65
1:A:437:THR:CG2	1:A:508:LEU:HD12	2.22	0.65
1:A:451:LEU:HD11	1:A:493:GLU:HG2	1.77	0.65
1:A:362:LEU:C	1:A:363:TYR:HD1	2.03	0.65
1:A:720:ASP:O	1:A:724:VAL:HG23	1.97	0.65
2:B:399:GLY:HA3	2:B:437:TRP:CE3	2.31	0.65
1:A:196:PHE:HB3	1:A:199:ILE:HG12	1.79	0.64
1:A:270:ILE:HG22	1:A:271:LYS:N	2.12	0.64
1:A:776:MET:HE3	1:A:776:MET:HA	1.79	0.64
1:A:460:GLN:O	1:A:464:GLU:HG3	1.97	0.63
1:A:392:LEU:HD12	1:A:415:VAL:HG23	1.79	0.63
1:A:671:TRP:HA	1:A:735:PHE:CE2	2.33	0.63
1:A:465:ALA:CB	1:A:479:LEU:HD23	2.28	0.63
1:A:662:VAL:HG13	1:A:748:VAL:HG22	1.80	0.63
1:A:319:THR:HB	1:A:572:SER:HB3	1.81	0.62
1:A:511:LEU:HD23	1:A:511:LEU:N	2.13	0.62
1:A:730:ILE:O	1:A:734:ILE:HD13	1.99	0.62
1:A:654:MET:HE1	1:A:776:MET:CG	2.26	0.62
1:A:761:TYR:CD1	1:A:809:ALA:HB1	2.35	0.62
1:A:363:TYR:CE2	1:A:734:ILE:HG23	2.35	0.61
2:B:307:LYS:O	2:B:308:ARG:HB2	2.00	0.61
2:B:406:SER:HB2	2:B:417:VAL:HG21	1.82	0.61
1:A:445:LEU:CB	2:B:359:LEU:HD23	2.30	0.61
1:A:353:LEU:HD13	1:A:565:LEU:HD23	1.83	0.61
1:A:654:MET:CE	1:A:776:MET:CG	2.79	0.61
2:B:324:VAL:HG23	2:B:331:ALA:HA	1.83	0.61
1:A:520:TYR:CE2	1:A:521:LEU:HD12	2.36	0.60
2:B:333:THR:HG22	2:B:334:VAL:N	2.16	0.60
1:A:346:SER:HB3	1:A:351:MET:CE	2.31	0.60
1:A:499:GLU:HG3	1:A:500:THR:N	2.16	0.59
1:A:537:GLU:OE2	1:A:544:LEU:HD13	2.03	0.59
1:A:568:ARG:NH1	1:A:699:LYS:HG3	2.17	0.59
1:A:716:GLU:HG2	1:A:750:ARG:HG2	1.84	0.59
1:A:346:SER:HA	1:A:351:MET:CE	2.33	0.59
1:A:231:PHE:CE1	1:A:249:VAL:HG12	2.36	0.59
2:B:307:LYS:HD2	2:B:308:ARG:HG3	1.83	0.59
2:B:307:LYS:HD2	2:B:308:ARG:NE	2.15	0.59
1:A:444:LEU:HD23	1:A:501:GLN:HB2	1.85	0.58
1:A:564:HIS:C	1:A:565:LEU:HD12	2.27	0.58
2:B:422:VAL:HA	2:B:425:ARG:HB2	1.84	0.58
1:A:363:TYR:HD2	1:A:734:ILE:CG1	2.14	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:455:ILE:HD11	1:A:490:LEU:O	2.04	0.58
1:A:470:PRO:O	1:A:472:ARG:N	2.37	0.58
1:A:633:GLN:HB2	1:A:634:PRO:HD3	1.85	0.58
1:A:188:MET:CE	1:A:200:ILE:HD12	2.23	0.57
1:A:695:TRP:HE3	1:A:697:LEU:HD21	1.69	0.57
2:B:311:PRO:HG2	2:B:314:MET:CG	2.24	0.57
1:A:485:ARG:HG3	1:A:485:ARG:NH1	2.19	0.57
1:A:724:VAL:HG11	1:A:746:THR:HG21	1.87	0.57
2:B:400:ARG:O	2:B:402:PHE:N	2.36	0.57
1:A:346:SER:CA	1:A:351:MET:HE3	2.34	0.56
2:B:340:MET:HA	2:B:343:VAL:HG22	1.85	0.56
1:A:418:LEU:HD11	2:B:321:VAL:HG22	1.86	0.56
1:A:421:LYS:NZ	2:B:320:ASP:OD2	2.38	0.56
2:B:395:ILE:HG22	2:B:433:VAL:HG12	1.88	0.56
1:A:188:MET:CE	1:A:200:ILE:HA	2.36	0.56
1:A:188:MET:HE1	1:A:200:ILE:N	2.20	0.56
1:A:380:GLN:O	1:A:384:ARG:HG3	2.05	0.56
1:A:402:ASN:O	1:A:403:ASN:HB2	2.06	0.55
1:A:183:LEU:HD23	1:A:189:THR:HG21	1.89	0.55
1:A:821:GLU:OE1	1:A:824:ARG:NH1	2.40	0.55
2:B:341:GLU:HG3	2:B:341:GLU:O	2.07	0.55
1:A:363:TYR:CD2	1:A:734:ILE:HG13	2.36	0.55
2:B:320:ASP:O	2:B:323:ALA:HB3	2.07	0.55
1:A:632:GLN:HG3	1:A:758:ARG:HD2	1.89	0.55
1:A:646:TRP:CZ3	1:A:647:LYS:HE2	2.42	0.55
1:A:667:ASP:CG	1:A:744:LYS:HE2	2.32	0.54
1:A:346:SER:HA	1:A:351:MET:HE3	1.89	0.54
1:A:374:LYS:NZ	1:A:525:ASP:OD1	2.36	0.54
1:A:356:ILE:HD11	1:A:566:THR:HG22	1.90	0.54
1:A:594:ARG:O	1:A:600:CYS:HB3	2.07	0.54
2:B:401:ASP:O	2:B:403:GLN:N	2.41	0.54
1:A:655:GLY:HA3	1:A:763:TYR:CZ	2.42	0.54
1:A:266:ILE:HD11	1:A:578:LEU:HD23	1.89	0.53
1:A:658:ASN:ND2	1:A:752:ARG:HB2	2.23	0.53
1:A:453:GLU:OE2	1:A:453:GLU:HA	2.07	0.53
1:A:356:ILE:HD11	1:A:566:THR:CG2	2.38	0.53
1:A:173:ALA:HB3	1:A:175:GLU:OE2	2.09	0.53
1:A:445:LEU:HB3	2:B:359:LEU:HD23	1.89	0.53
1:A:195:CYS:C	1:A:197:PRO:HD3	2.34	0.53
1:A:282:ILE:HG21	1:A:602:VAL:HG21	1.91	0.53
1:A:501:GLN:O	1:A:501:GLN:HG2	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:315:PHE:O	2:B:316:LEU:HD23	2.09	0.53
2:B:348:GLN:O	2:B:348:GLN:HG2	2.09	0.52
1:A:572:SER:O	1:A:575:PRO:HD2	2.08	0.52
2:B:388:GLN:HB3	2:B:428:PHE:HE2	1.74	0.52
1:A:449:VAL:HG23	2:B:363:LEU:CD2	2.39	0.52
1:A:487:LEU:HD23	2:B:372:LEU:HD11	1.92	0.52
2:B:307:LYS:CD	2:B:308:ARG:HE	2.20	0.52
1:A:183:LEU:CD2	1:A:189:THR:HG21	2.40	0.52
1:A:363:TYR:CD2	1:A:734:ILE:HG23	2.45	0.51
1:A:672:ASP:O	1:A:675:VAL:HG22	2.09	0.51
1:A:346:SER:CA	1:A:351:MET:CE	2.88	0.51
1:A:401:LEU:O	1:A:402:ASN:HB2	2.10	0.51
1:A:538:PHE:HB2	1:A:708:ALA:HB2	1.92	0.51
1:A:275:THR:HG23	1:A:276:LYS:H	1.76	0.51
1:A:180:GLN:HA	1:A:180:GLN:OE1	2.10	0.51
1:A:456:LYS:HA	2:B:370:TYR:HE2	1.75	0.51
1:A:488:THR:HA	1:A:491:CYS:HB2	1.92	0.51
1:A:601:GLU:HA	1:A:616:TYR:O	2.10	0.50
1:A:231:PHE:CZ	1:A:250:HIS:HB2	2.47	0.50
1:A:363:TYR:CD2	1:A:734:ILE:CG1	2.93	0.50
2:B:432:GLU:O	2:B:435:GLN:HG2	2.11	0.50
1:A:418:LEU:HD22	2:B:320:ASP:HB3	1.94	0.50
1:A:513:ALA:C	1:A:515:PRO:HD3	2.37	0.50
1:A:428:ILE:O	1:A:432:LYS:HB2	2.11	0.50
2:B:415:VAL:HG13	2:B:416:GLN:N	2.24	0.50
1:A:393:SER:HB2	1:A:549:LEU:HD21	1.92	0.50
1:A:538:PHE:CE1	1:A:706:LEU:HD13	2.47	0.50
1:A:364:GLU:OE2	1:A:524:ARG:NH1	2.46	0.49
1:A:695:TRP:CE3	1:A:697:LEU:HD11	2.47	0.49
1:A:230:THR:HG21	1:A:270:ILE:HG21	1.94	0.49
2:B:369:PRO:HG2	2:B:370:TYR:CE1	2.47	0.49
1:A:470:PRO:CD	1:A:471:PRO:HD2	2.42	0.49
1:A:572:SER:C	1:A:575:PRO:HD2	2.38	0.49
1:A:548:SER:O	1:A:552:TRP:HB3	2.12	0.49
2:B:436:GLU:HA	2:B:436:GLU:OE2	2.12	0.49
1:A:319:THR:HG22	1:A:321:ARG:HG3	1.95	0.49
1:A:372:LYS:O	1:A:376:GLU:HG3	2.12	0.49
1:A:786:ILE:CG2	1:A:787:PRO:HD2	2.42	0.49
1:A:781:THR:HG22	1:A:781:THR:O	2.12	0.48
1:A:438:GLN:CG	1:A:508:LEU:HD11	2.40	0.48
1:A:445:LEU:HB2	2:B:359:LEU:HD23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:MET:HE1	1:A:200:ILE:CA	2.43	0.48
1:A:677:LEU:HB2	1:A:693:LEU:HD11	1.95	0.48
1:A:456:LYS:CG	2:B:370:TYR:HE2	2.23	0.48
1:A:340:ASN:OD1	1:A:342:MET:N	2.41	0.47
1:A:325:TYR:O	1:A:326:VAL:HG23	2.14	0.47
1:A:363:TYR:HD1	1:A:363:TYR:N	2.12	0.47
1:A:487:LEU:HD23	2:B:372:LEU:CD1	2.44	0.47
1:A:510:GLU:HG2	1:A:511:LEU:N	2.28	0.47
1:A:736:GLY:C	1:A:738:SER:H	2.23	0.47
1:A:346:SER:CB	1:A:351:MET:CE	2.87	0.47
2:B:411:ASN:C	2:B:412:LYS:HD2	2.38	0.47
1:A:563:SER:O	1:A:565:LEU:HD13	2.15	0.47
1:A:568:ARG:HH12	1:A:699:LYS:N	2.12	0.47
1:A:180:GLN:OE1	1:A:339:GLY:N	2.43	0.47
1:A:196:PHE:HB3	1:A:199:ILE:CG1	2.44	0.47
1:A:452:LYS:HD2	2:B:366:GLY:O	2.14	0.47
1:A:363:TYR:N	1:A:363:TYR:CD1	2.83	0.47
1:A:392:LEU:HD22	1:A:398:PHE:CB	2.45	0.47
1:A:537:GLU:HG2	1:A:544:LEU:HD11	1.97	0.47
1:A:537:GLU:CG	1:A:544:LEU:HD13	2.41	0.47
1:A:209:VAL:O	1:A:213:ILE:HG13	2.15	0.46
1:A:458:LEU:HD23	1:A:458:LEU:HA	1.61	0.46
1:A:671:TRP:O	1:A:673:PRO:CD	2.59	0.46
1:A:677:LEU:HA	1:A:694:PHE:O	2.16	0.46
1:A:317:VAL:HG13	1:A:571:TYR:HB3	1.97	0.46
2:B:413:SER:O	2:B:415:VAL:N	2.49	0.46
1:A:362:LEU:C	1:A:363:TYR:CD1	2.90	0.46
1:A:411:ALA:CB	1:A:549:LEU:HD13	2.45	0.46
1:A:458:LEU:HD11	1:A:486:ASP:HB3	1.97	0.46
1:A:465:ALA:HB2	1:A:479:LEU:CD2	2.45	0.46
1:A:460:GLN:OE1	1:A:460:GLN:HA	2.15	0.46
1:A:667:ASP:OD2	1:A:744:LYS:CE	2.64	0.46
1:A:574:VAL:HB	1:A:575:PRO:HD3	1.98	0.46
1:A:807:TYR:N	1:A:808:PRO:HD3	2.31	0.46
2:B:400:ARG:O	2:B:402:PHE:CD1	2.68	0.46
1:A:245:ASP:OD1	1:A:247:VAL:CG1	2.64	0.45
1:A:538:PHE:CE1	1:A:706:LEU:CD1	2.99	0.45
2:B:334:VAL:O	2:B:338:LEU:HG	2.16	0.45
2:B:388:GLN:O	2:B:391:ALA:HB3	2.17	0.45
1:A:568:ARG:HG2	1:A:568:ARG:HH11	1.81	0.45
1:A:728:LEU:HA	1:A:731:LEU:HD12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:ARG:HG2	1:A:326:VAL:HG22	1.98	0.45
2:B:388:GLN:O	2:B:392:VAL:HG12	2.16	0.45
1:A:808:PRO:O	1:A:810:THR:HG23	2.16	0.45
1:A:327:ALA:CB	1:A:663:VAL:HG11	2.47	0.45
1:A:378:VAL:HG11	1:A:528:ILE:HG22	1.98	0.45
2:B:307:LYS:O	2:B:308:ARG:CB	2.64	0.45
2:B:324:VAL:HG22	2:B:324:VAL:O	2.17	0.45
1:A:776:MET:HB3	1:A:803:THR:HG22	1.99	0.45
2:B:331:ALA:C	2:B:333:THR:H	2.25	0.44
2:B:424:TYR:HA	2:B:427:ARG:NH1	2.33	0.44
1:A:455:ILE:HG23	1:A:487:LEU:HD11	1.99	0.44
1:A:478:PHE:O	1:A:482:SER:HB3	2.17	0.44
1:A:660:ASN:OD1	1:A:749:SER:OG	2.36	0.44
2:B:402:PHE:CE2	2:B:418:LYS:HG3	2.53	0.44
1:A:469:LYS:C	1:A:471:PRO:HD2	2.43	0.44
1:A:667:ASP:OD2	1:A:744:LYS:HE2	2.17	0.44
2:B:424:TYR:O	2:B:426:ARG:N	2.51	0.44
1:A:501:GLN:O	1:A:505:GLU:HG3	2.16	0.44
1:A:583:ASP:OD1	1:A:585:LYS:NZ	2.50	0.44
1:A:213:ILE:CD1	1:A:248:LEU:HD23	2.48	0.44
2:B:368:GLU:OE2	2:B:371:ARG:NH2	2.51	0.44
2:B:388:GLN:HB3	2:B:428:PHE:CE2	2.53	0.44
1:A:384:ARG:NH2	2:B:312:LYS:O	2.44	0.43
1:A:760:SER:O	3:A:901:FAD:HM81	2.18	0.43
2:B:314:MET:HE3	2:B:314:MET:HB3	1.71	0.43
1:A:691:LEU:CD2	1:A:705:ALA:HB1	2.48	0.43
2:B:337:GLN:HG3	2:B:338:LEU:HD23	2.00	0.43
1:A:188:MET:CE	1:A:200:ILE:CA	2.96	0.43
1:A:230:THR:HG23	1:A:270:ILE:HD12	1.99	0.43
1:A:633:GLN:HB2	1:A:634:PRO:CD	2.47	0.43
2:B:324:VAL:HG23	2:B:331:ALA:CA	2.48	0.43
2:B:383:TRP:CZ2	2:B:420:PHE:HB2	2.54	0.43
1:A:543:PRO:HG3	1:A:710:GLU:HG2	2.01	0.43
1:A:238:LEU:HD23	1:A:238:LEU:HA	1.86	0.43
1:A:490:LEU:N	1:A:490:LEU:HD23	2.33	0.43
1:A:238:LEU:HD12	1:A:249:VAL:HG21	2.00	0.43
1:A:456:LYS:HA	2:B:370:TYR:CE2	2.53	0.43
1:A:534:ALA:HA	1:A:537:GLU:HG3	2.01	0.43
2:B:424:TYR:CE1	2:B:427:ARG:NH2	2.87	0.43
1:A:474:ILE:HD12	1:A:474:ILE:HA	1.90	0.43
1:A:494:TYR:CD1	1:A:494:TYR:C	2.97	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:541:ALA:O	1:A:657:GLY:HA3	2.18	0.43
1:A:786:ILE:HG23	1:A:787:PRO:HD2	2.00	0.43
1:A:412:LEU:HA	1:A:412:LEU:HD23	1.83	0.43
1:A:718:ILE:HG22	1:A:723:ILE:HG13	2.01	0.43
1:A:242:TYR:C	1:A:244:SER:H	2.27	0.42
1:A:732:LYS:C	1:A:734:ILE:H	2.25	0.42
1:A:760:SER:CA	3:A:901:FAD:HM81	2.49	0.42
1:A:714:ILE:O	1:A:716:GLU:N	2.52	0.42
1:A:392:LEU:HD22	1:A:398:PHE:HB3	2.00	0.42
2:B:331:ALA:O	2:B:333:THR:N	2.53	0.42
1:A:306:LEU:HD11	1:A:582:LEU:HD13	2.02	0.42
1:A:456:LYS:CG	2:B:370:TYR:CE2	2.98	0.42
1:A:575:PRO:O	1:A:576:VAL:C	2.63	0.42
1:A:691:LEU:HD22	1:A:705:ALA:HB1	2.01	0.42
1:A:434:ILE:HG22	2:B:349:ILE:CD1	2.49	0.42
1:A:694:PHE:HA	1:A:704:LEU:O	2.19	0.42
1:A:695:TRP:CE3	1:A:697:LEU:HD21	2.53	0.42
2:B:383:TRP:CD2	2:B:412:LYS:NZ	2.88	0.42
1:A:188:MET:HB2	1:A:200:ILE:CD1	2.50	0.42
1:A:670:PHE:CD1	1:A:670:PHE:C	2.98	0.42
2:B:438:GLU:O	2:B:439:ALA:C	2.61	0.42
1:A:367:GLY:C	1:A:368:GLN:CD	2.88	0.41
2:B:368:GLU:C	2:B:370:TYR:H	2.28	0.41
2:B:390:LEU:HD23	2:B:390:LEU:HA	1.58	0.41
1:A:196:PHE:N	1:A:197:PRO:HD3	2.34	0.41
1:A:210:PHE:CE1	1:A:252:VAL:HG22	2.55	0.41
1:A:266:ILE:CD1	1:A:578:LEU:HD23	2.50	0.41
1:A:386:LEU:HD23	1:A:386:LEU:HA	1.90	0.41
1:A:695:TRP:HB2	1:A:704:LEU:HB2	2.02	0.41
1:A:451:LEU:HA	1:A:451:LEU:HD23	1.73	0.41
2:B:338:LEU:HD23	2:B:338:LEU:N	2.35	0.41
2:B:418:LYS:O	2:B:421:PHE:HB2	2.20	0.41
1:A:191:GLN:HG2	1:A:255:TYR:OH	2.20	0.41
1:A:780:ILE:HB	1:A:796:LEU:HB3	2.02	0.41
1:A:786:ILE:H	1:A:786:ILE:CD1	2.29	0.41
1:A:801:GLU:HG3	1:A:809:ALA:HA	2.01	0.41
1:A:493:GLU:O	1:A:496:GLU:HG2	2.21	0.41
2:B:307:LYS:HD2	2:B:308:ARG:CG	2.51	0.41
1:A:367:GLY:N	1:A:368:GLN:NE2	2.69	0.41
1:A:565:LEU:N	1:A:565:LEU:CD1	2.83	0.41
1:A:622:LEU:HD11	1:A:821:GLU:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:ASP:CG	1:A:251:ARG:HH22	2.25	0.41
1:A:295:ARG:NH2	1:A:580:GLU:O	2.54	0.41
1:A:489:ALA:C	1:A:491:CYS:N	2.78	0.41
1:A:755:PRO:HA	1:A:758:ARG:HE	1.86	0.41
2:B:402:PHE:CD2	2:B:418:LYS:HG3	2.56	0.41
1:A:520:TYR:CE2	1:A:521:LEU:CD1	3.04	0.41
1:A:760:SER:HA	3:A:901:FAD:HM81	2.03	0.41
1:A:420:GLU:OE2	1:A:526:ARG:NH1	2.50	0.40
1:A:356:ILE:CG1	1:A:566:THR:HG23	2.51	0.40
1:A:725:GLY:O	1:A:726:ARG:C	2.64	0.40
1:A:485:ARG:C	1:A:485:ARG:HD2	2.46	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	658/660 (100%)	600 (91%)	47 (7%)	11 (2%)	7	15
2	B	132/134 (98%)	100 (76%)	20 (15%)	12 (9%)	0	0
All	All	790/794 (100%)	700 (89%)	67 (8%)	23 (3%)	3	8

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	240	ALA
1	A	350	ASN
2	B	308	ARG
2	B	319	GLU
2	B	331	ALA
2	B	375	VAL
2	B	429	ASN

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Mol	Chain	Res	Type
1	A	490	LEU
1	A	792	PRO
2	B	318	GLN
2	B	332	THR
2	B	377	GLN
2	B	402	PHE
2	B	425	ARG
1	A	471	PRO
1	A	690	GLU
1	A	729	ALA
1	A	365	ALA
1	A	516	PRO
1	A	715	MET
1	A	737	SER
2	B	401	ASP
2	B	415	VAL

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	558/560 (100%)	529 (95%)	29 (5%)	21	43
2	B	112/118 (95%)	106 (95%)	6 (5%)	20	42
All	All	670/678 (99%)	635 (95%)	35 (5%)	21	43

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	191	GLN
1	A	198	ASP
1	A	232	GLU
1	A	244	SER
1	A	326	VAL
1	A	359	LYS
1	A	392	LEU

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Mol	Chain	Res	Type
1	A	443	GLU
1	A	446	ASN
1	A	449	VAL
1	A	491	CYS
1	A	509	GLN
1	A	538	PHE
1	A	563	SER
1	A	565	LEU
1	A	598	SER
1	A	611	SER
1	A	624	THR
1	A	662	VAL
1	A	663	VAL
1	A	667	ASP
1	A	684	THR
1	A	685	THR
1	A	727	CYS
1	A	738	SER
1	A	771	ASN
1	A	781	THR
1	A	793	ILE
1	A	815	LEU
2	B	333	THR
2	B	371	ARG
2	B	375	VAL
2	B	385	THR
2	B	422	VAL
2	B	426	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	191	GLN
1	A	237	GLN
1	A	368	GLN
1	A	380	GLN
1	A	430	HIS
1	A	438	GLN
1	A	446	ASN
1	A	612	GLN
1	A	652	GLN
2	B	435	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 ⓘ

3 ligands are 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	$\# Z > 2$
3	FAD	A	901	-	58,58,58	0.47	0	85,89,89	0.72	2 (2%)

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	FAD	A	901	-	-	7/34/50/50	0/6/6/6

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	FAD	O5'-P-O1P	-3.64	94.51	108.94
3	A	901	FAD	O2A-PA-O1A	2.08	122.11	112.44

There are no chirality outliers.

All (7) torsion outliers are listed below:

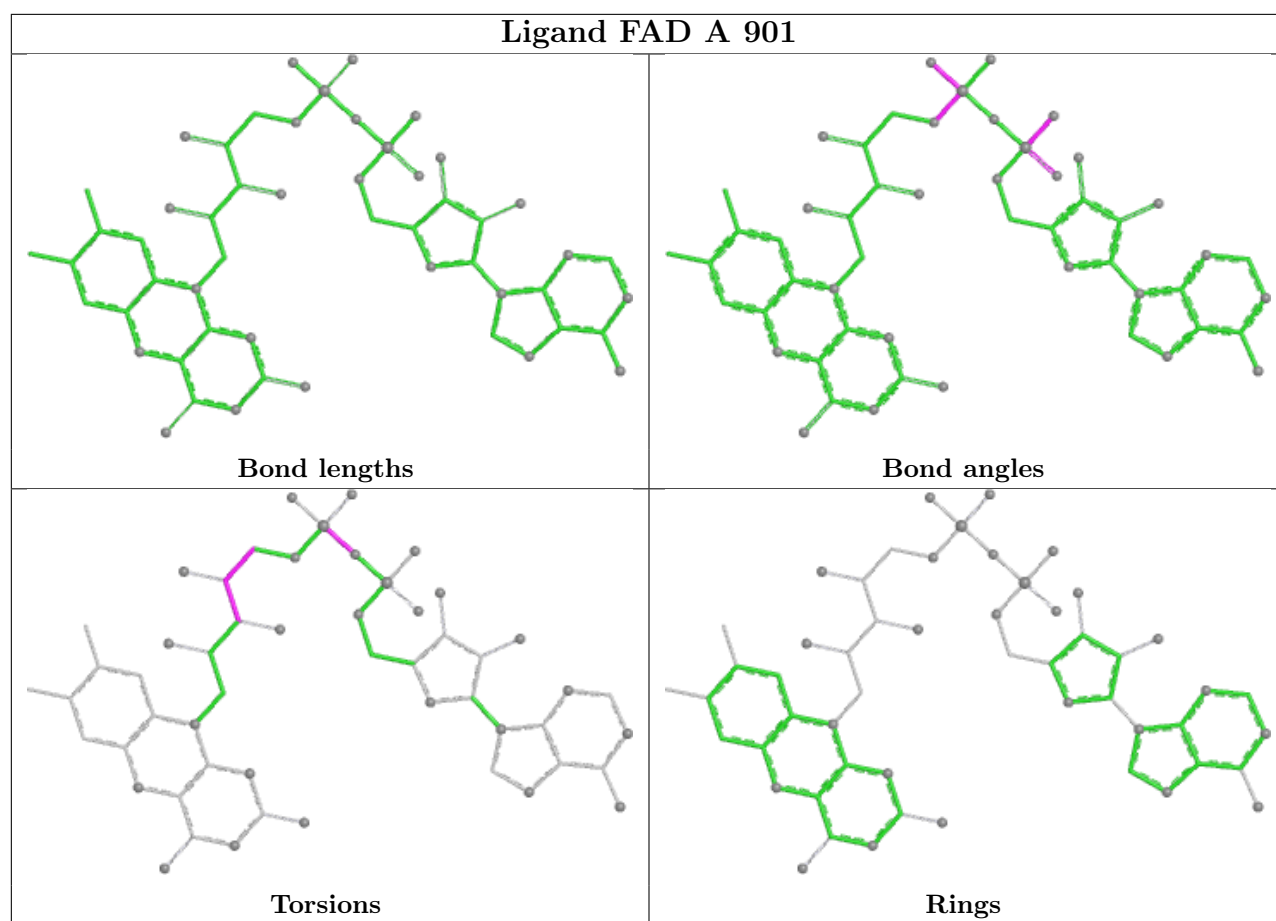
Mol	Chain	Res	Type	Atoms
3	A	901	FAD	PA-O3P-P-O5'
3	A	901	FAD	C2'-C3'-C4'-C5'
3	A	901	FAD	C2'-C3'-C4'-O4'
3	A	901	FAD	C3'-C4'-C5'-O5'
3	A	901	FAD	O4'-C4'-C5'-O5'
3	A	901	FAD	O3'-C3'-C4'-C5'
3	A	901	FAD	O3'-C3'-C4'-O4'

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	FAD	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	660/660 (100%)	0.28	38 (5%) 29 22	38, 67, 98, 116	0
2	B	134/134 (100%)	1.11	24 (17%) 3 3	70, 97, 113, 129	0
All	All	794/794 (100%)	0.42	62 (7%) 19 14	38, 74, 105, 129	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	365	ALA	7.8
2	B	375	VAL	5.9
2	B	377	GLN	5.1
2	B	307	LYS	4.9
1	A	369	ALA	4.4
2	B	374	GLU	4.3
1	A	368	GLN	4.1
1	A	366	ASN	4.0
1	A	173	ALA	3.8
1	A	275	THR	3.8
1	A	667	ASP	3.7
2	B	332	THR	3.7
1	A	633	GLN	3.6
2	B	440	GLU	3.4
1	A	359	LYS	3.4
1	A	276	LYS	3.4
1	A	269	ARG	3.3
2	B	382	ARG	3.3
1	A	373	GLU	3.1
2	B	378	LYS	3.0
1	A	492	LYS	3.0
1	A	527	GLN	3.0
1	A	380	GLN	2.9
1	A	791	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	175	GLU	2.8
2	B	439	ALA	2.7
1	A	471	PRO	2.7
1	A	610	THR	2.7
2	B	376	ILE	2.7
1	A	470	PRO	2.7
2	B	379	CYS	2.6
1	A	832	ALA	2.6
1	A	239	GLU	2.5
2	B	308	ARG	2.5
1	A	367	GLY	2.5
1	A	668	ARG	2.5
1	A	377	MET	2.5
1	A	358	GLN	2.5
1	A	699	LYS	2.4
2	B	329	THR	2.4
2	B	418	LYS	2.4
2	B	340	MET	2.4
1	A	508	LEU	2.4
1	A	439	GLU	2.4
1	A	512	GLU	2.4
2	B	423	ASN	2.3
2	B	312	LYS	2.3
2	B	422	VAL	2.3
1	A	682	GLY	2.3
1	A	453	GLU	2.3
1	A	236	GLN	2.2
2	B	380	ASN	2.2
1	A	324	ASN	2.2
2	B	383	TRP	2.1
1	A	634	PRO	2.1
1	A	518	ASP	2.1
2	B	399	GLY	2.1
1	A	524	ARG	2.1
2	B	381	ALA	2.0
2	B	431	ASP	2.0
1	A	205	GLN	2.0
2	B	438	GLU	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 [i](#)

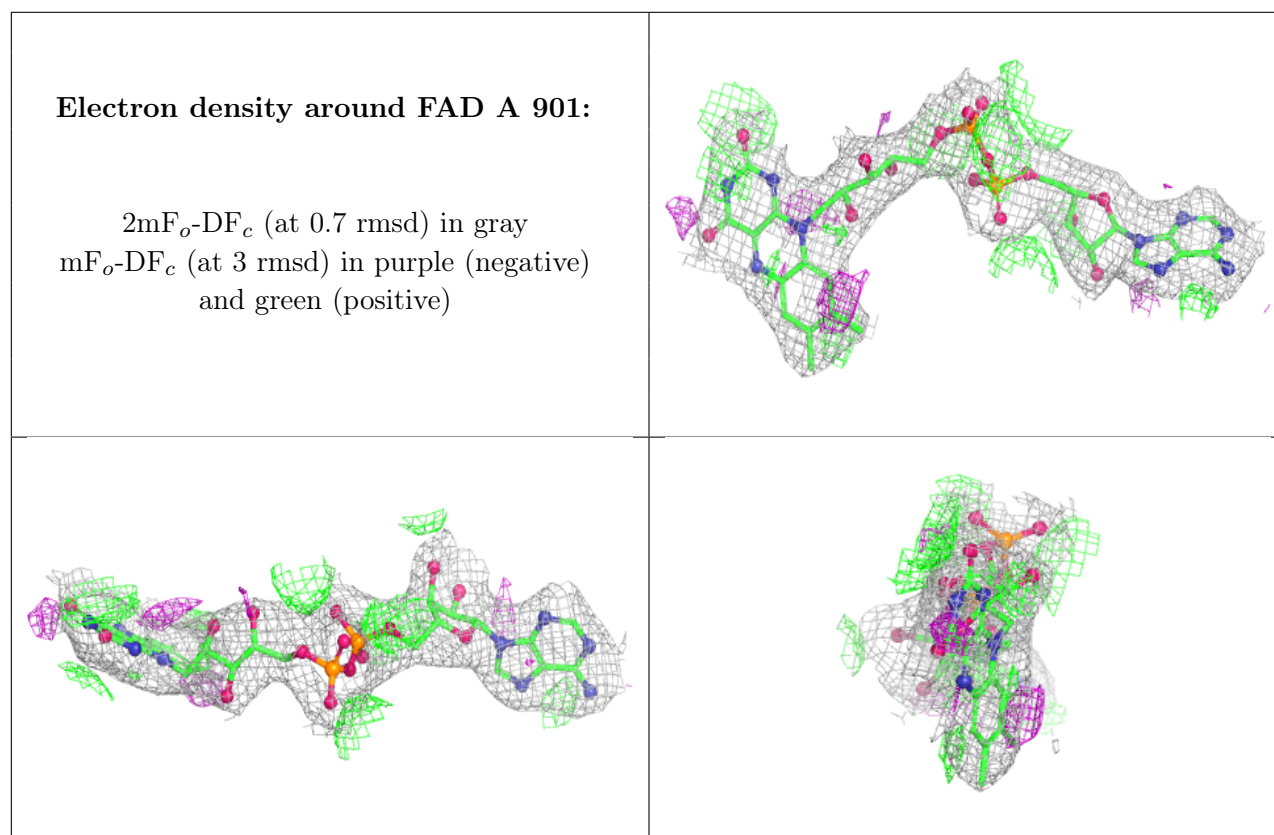
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
4	A1LZJ	A	902[A]	33/33	0.95	0.10	44,57,73,78	1
4	A1LZJ	A	902[B]	33/33	0.95	0.10	44,57,73,78	1
3	FAD	A	901	53/53	0.97	0.08	42,48,60,65	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.