



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:57 AM UTC

PDB ID : 8FSK / pdb_00008fsk
Title : LSD1-CoREST in complex with T18, short soaking
Authors : Caroli, J.; Mattevi, A.
Deposited on : 2023-01-10
Resolution : 3.13 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

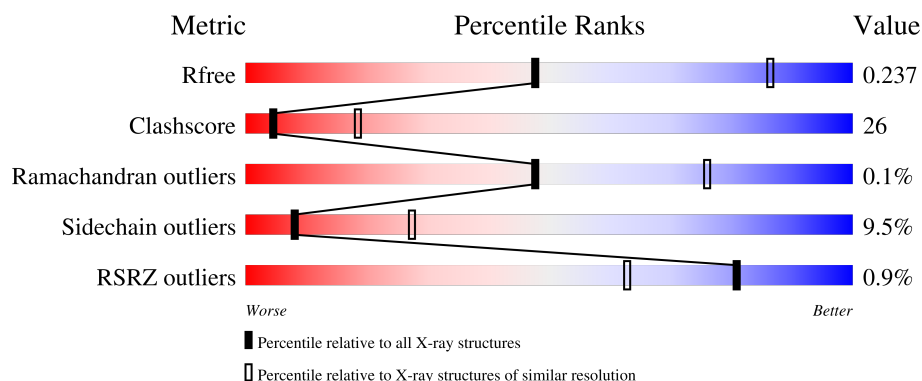
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.13 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	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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	871	50% 22% • • 24%
2	B	144	40% 40% 10% • 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6365 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 Lysine-specific histone demethylase 1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	666	Total	C	N	O	S	0	0	0
			5217	3324	906	967	20			

There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	GLY	-	expression tag	UNP O60341
A	-17	SER	-	expression tag	UNP O60341
A	-16	SER	-	expression tag	UNP O60341
A	-15	HIS	-	expression tag	UNP O60341
A	-14	HIS	-	expression tag	UNP O60341
A	-13	HIS	-	expression tag	UNP O60341
A	-12	HIS	-	expression tag	UNP O60341
A	-11	HIS	-	expression tag	UNP O60341
A	-10	HIS	-	expression tag	UNP O60341
A	-9	SER	-	expression tag	UNP O60341
A	-8	SER	-	expression tag	UNP O60341
A	-7	GLY	-	expression tag	UNP O60341
A	-6	LEU	-	expression tag	UNP O60341
A	-5	VAL	-	expression tag	UNP O60341
A	-4	PRO	-	expression tag	UNP O60341
A	-3	ARG	-	expression tag	UNP O60341
A	-2	GLY	-	expression tag	UNP O60341
A	-1	SER	-	expression tag	UNP O60341
A	0	HIS	-	expression tag	UNP O60341

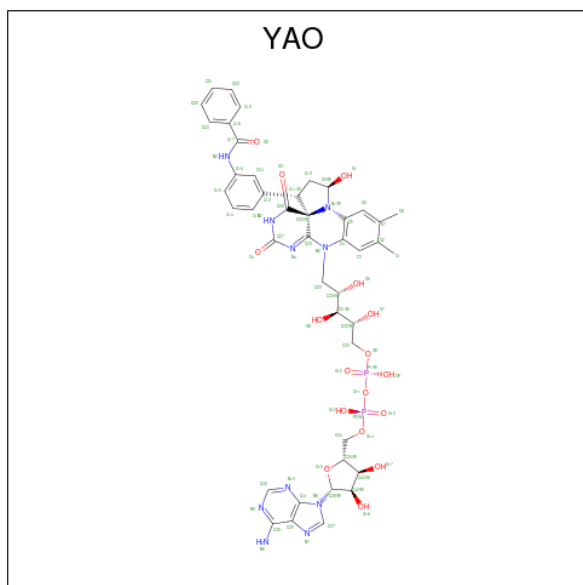
- Molecule 2 is a protein called REST corepressor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	133	Total	C	N	O	S	0	0	0
			1076	676	194	203	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	297	GLY	-	expression tag	UNP Q9UKL0
B	298	PRO	-	expression tag	UNP Q9UKL0
B	299	LEU	-	expression tag	UNP Q9UKL0
B	300	GLY	-	expression tag	UNP Q9UKL0
B	301	SER	-	expression tag	UNP Q9UKL0
B	302	PRO	-	expression tag	UNP Q9UKL0
B	303	GLU	-	expression tag	UNP Q9UKL0
B	304	PHE	-	expression tag	UNP Q9UKL0

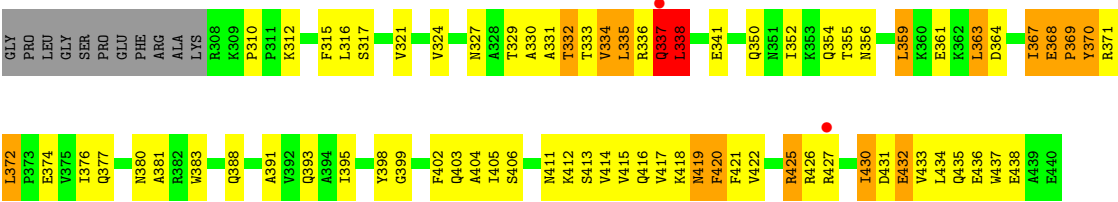
- Molecule 3 is [(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-3,4-dihydroxyoxolan-2-yl]methyl (2R,3S,4S)-5-[(1R,3S,3aS,13R)-3-(3-benzamidophenyl)-1-hydroxy-10,11-dimethyl-4,6-dioxo-2,3,5,6-tetrahydro-1H-benzo[g]pyrrolo[2,1-e]pteridin-8(4H)-yl]-2,3,4-trihydroxypentyl dihydrogen diphosphate (non-preferred name) (CCD ID: YAO) (formula: C₄₃H₄₈N₁₀O₁₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			72	43	10	17	2		

- Molecule 1: Lysine-specific histone demethylase 1A





4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	119.39Å 179.91Å 235.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.70 – 3.13 59.70 – 3.13	Depositor EDS
% Data completeness (in resolution range)	98.9 (59.70-3.13) 98.9 (59.70-3.13)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 3.13Å)	Xtriage
Refinement program	PHENIX 1.18.2	Depositor
R, R_{free}	0.199 , 0.223 0.218 , 0.237	Depositor DCC
R_{free} test set	2000 reflections (4.43%)	wwPDB-VP
Wilson B-factor (Å ²)	116.1	Xtriage
Anisotropy	0.389	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 86.5	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	0.95	EDS
Total number of atoms	6365	wwPDB-VP
Average B, all atoms (Å ²)	110.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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: YAO

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	0.94	20/5331 (0.4%)	1.17	20/7232 (0.3%)
2	B	1.08	3/1091 (0.3%)	1.45	12/1471 (0.8%)
All	All	0.96	23/6422 (0.4%)	1.22	32/8703 (0.4%)

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	659	LEU	C-O	-8.58	1.14	1.23
2	B	369	PRO	C-O	-8.34	1.14	1.24
1	A	479	LEU	C-O	-7.05	1.15	1.24
1	A	660	ASN	C-O	-6.85	1.14	1.23
1	A	201	SER	CA-CB	-6.60	1.42	1.53
1	A	326	VAL	C-O	-6.35	1.16	1.24
1	A	200	ILE	C-O	-6.18	1.16	1.24
1	A	278	THR	C-O	-5.89	1.17	1.24
1	A	192	GLU	C-O	-5.82	1.17	1.24
1	A	662	VAL	C-O	-5.72	1.18	1.24
1	A	193	ALA	C-O	-5.67	1.17	1.24
1	A	470	PRO	C-O	-5.65	1.20	1.24
1	A	661	LYS	C-O	-5.61	1.17	1.23
1	A	194	ALA	C-O	-5.58	1.17	1.24
1	A	320	PHE	C-O	-5.55	1.17	1.24
1	A	808	PRO	C-O	-5.53	1.17	1.23
1	A	504	LEU	C-O	-5.40	1.17	1.24
1	A	256	LEU	C-O	-5.39	1.17	1.24
1	A	312	ARG	C-O	-5.35	1.17	1.23
1	A	480	VAL	C-O	-5.19	1.18	1.24
1	A	191	GLN	C-O	-5.12	1.18	1.24
2	B	406	SER	C-O	-5.11	1.18	1.24
2	B	413	SER	CA-CB	-5.05	1.45	1.53

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	421	PHE	CB-CA-C	-8.85	96.98	110.88
2	B	431	ASP	N-CA-C	-7.96	102.19	111.03
2	B	337	GLN	CB-CA-C	-7.65	98.87	110.88
1	A	661	LYS	CB-CA-C	-7.11	93.78	109.56
1	A	270	ILE	N-CA-C	-6.97	104.39	110.74
2	B	364	ASP	CA-CB-CG	-6.96	105.64	112.60
2	B	369	PRO	CB-CA-C	-6.67	102.37	112.11
1	A	200	ILE	CB-CA-C	-6.64	101.86	112.16
1	A	381	GLU	CA-C-O	-6.43	114.07	120.82
1	A	635	PRO	CB-CA-C	-6.30	103.70	111.64
1	A	324	ASN	N-CA-C	-6.22	105.34	113.17
2	B	329	THR	CB-CA-C	-6.19	98.86	110.01
1	A	279	GLY	CA-C-O	-6.06	117.38	122.29
2	B	417	VAL	CB-CA-C	-5.94	104.12	112.14
2	B	333	THR	CB-CA-C	-5.87	101.67	110.88
2	B	338	LEU	N-CA-C	-5.73	105.03	111.28
2	B	332	THR	CB-CA-C	-5.70	101.32	110.79
1	A	381	GLU	O-C-N	5.61	127.85	122.07
2	B	426	ARG	O-C-N	5.59	128.13	122.09
1	A	349	VAL	CA-C-O	-5.57	114.86	120.31
1	A	470	PRO	CB-CA-C	-5.54	105.97	111.17
1	A	546	THR	CA-CB-OG1	-5.51	101.34	109.60
1	A	803	THR	CA-CB-OG1	-5.43	101.45	109.60
1	A	274	PRO	CB-CA-C	-5.34	104.57	111.46
1	A	278	THR	CB-CA-C	-5.29	100.02	109.71
1	A	642	PRO	CB-CA-C	-5.26	104.49	111.23
1	A	485	ARG	CB-CA-C	-5.26	102.62	110.88
1	A	316	ARG	CA-C-N	-5.16	116.80	121.65
1	A	316	ARG	C-N-CA	-5.16	116.80	121.65
2	B	432	GLU	CA-C-O	-5.04	115.50	120.90
1	A	274	PRO	CA-C-N	5.03	127.01	120.28
1	A	274	PRO	C-N-CA	5.03	127.01	120.28

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	5217	0	5252	277	0
2	B	1076	0	1091	87	0
3	A	72	0	0	4	0
All	All	6365	0	6343	332	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 (332) 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:193:ALA:CB	1:A:200:ILE:CD1	2.02	1.37
1:A:193:ALA:HB1	1:A:200:ILE:CD1	1.59	1.30
1:A:193:ALA:CA	1:A:200:ILE:HD11	1.65	1.26
1:A:438:GLN:OE1	1:A:508:LEU:HD12	1.48	1.13
1:A:456:LYS:HA	2:B:370:TYR:CE1	1.85	1.09
1:A:438:GLN:OE1	1:A:508:LEU:CD1	2.01	1.08
1:A:231:PHE:HE1	1:A:249:VAL:HG12	1.02	1.07
1:A:437:THR:CG2	1:A:508:LEU:HD21	1.84	1.07
1:A:793:ILE:HG23	1:A:828:GLN:NE2	1.69	1.06
1:A:231:PHE:CE1	1:A:249:VAL:HG12	1.92	1.05
1:A:193:ALA:CB	1:A:200:ILE:HD13	1.80	1.05
1:A:193:ALA:CB	1:A:200:ILE:HD11	1.74	1.04
1:A:192:GLU:OE2	1:A:214:ARG:NH1	1.91	1.04
1:A:193:ALA:HA	1:A:200:ILE:HD11	1.35	1.03
1:A:656:PHE:CE2	1:A:759:GLY:HA3	1.95	1.02
1:A:319:THR:HB	1:A:572:SER:HB3	1.37	1.01
1:A:193:ALA:CA	1:A:200:ILE:CD1	2.35	0.99
1:A:437:THR:HG22	1:A:508:LEU:HD21	1.41	0.98
1:A:193:ALA:HB2	1:A:200:ILE:HD13	1.41	0.98
1:A:667:ASP:CG	1:A:744:LYS:NZ	2.23	0.97
1:A:469:LYS:HE2	1:A:469:LYS:HA	1.47	0.96
1:A:793:ILE:HD12	1:A:793:ILE:H	1.26	0.95
1:A:193:ALA:HA	1:A:200:ILE:CD1	1.94	0.95
1:A:231:PHE:HE1	1:A:249:VAL:CG1	1.80	0.94
1:A:456:LYS:HA	2:B:370:TYR:HE1	1.30	0.94
1:A:193:ALA:HB1	1:A:200:ILE:HD12	1.49	0.92
1:A:801:GLU:HG3	1:A:809:ALA:HA	1.50	0.91
1:A:265:GLY:O	1:A:295:ARG:HD3	1.70	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:399:GLY:N	2:B:437:TRP:NE1	2.19	0.89
1:A:456:LYS:CA	2:B:370:TYR:HE1	1.87	0.88
1:A:568:ARG:NH1	1:A:699:LYS:HG2	1.88	0.87
2:B:425:ARG:HG3	2:B:425:ARG:HH11	1.38	0.86
1:A:384:ARG:NH2	2:B:312:LYS:O	2.07	0.86
1:A:566:THR:HG21	1:A:697:LEU:HD13	1.57	0.85
1:A:793:ILE:CG2	1:A:828:GLN:NE2	2.38	0.85
1:A:437:THR:CG2	1:A:508:LEU:CD2	2.54	0.85
1:A:695:TRP:HE1	1:A:706:LEU:HD13	1.40	0.85
1:A:568:ARG:CZ	1:A:699:LYS:HG2	2.07	0.84
1:A:350:ASN:O	1:A:350:ASN:ND2	2.11	0.82
1:A:353:LEU:HD13	1:A:565:LEU:HD22	1.61	0.81
1:A:667:ASP:CG	1:A:744:LYS:HZ1	1.85	0.81
1:A:456:LYS:CA	2:B:370:TYR:CE1	2.63	0.81
1:A:193:ALA:HB1	1:A:200:ILE:HD11	1.41	0.81
1:A:319:THR:CB	1:A:572:SER:HB3	2.10	0.81
1:A:456:LYS:CB	2:B:370:TYR:HE1	1.93	0.80
1:A:750:ARG:HG3	1:A:750:ARG:HH11	1.46	0.80
1:A:456:LYS:HA	2:B:370:TYR:CD1	2.17	0.79
1:A:320:PHE:CE1	1:A:747:VAL:HG21	2.18	0.78
1:A:392:LEU:HD11	2:B:316:LEU:HD22	1.63	0.78
1:A:695:TRP:HE1	1:A:706:LEU:CD1	1.96	0.78
1:A:667:ASP:HB3	1:A:744:LYS:NZ	1.99	0.77
2:B:359:LEU:HD23	2:B:359:LEU:N	1.99	0.77
1:A:647:LYS:NZ	1:A:776:MET:O	2.19	0.76
1:A:342:MET:HE2	1:A:571:TYR:OH	1.86	0.76
1:A:351:MET:HA	1:A:569:ASN:HD21	1.51	0.75
1:A:569:ASN:OD1	1:A:569:ASN:N	2.18	0.75
1:A:308:GLU:OE1	3:A:901:YAO:O17	2.04	0.75
1:A:815:LEU:HD12	1:A:815:LEU:O	1.86	0.75
2:B:335:LEU:N	2:B:335:LEU:HD23	2.01	0.75
1:A:312:ARG:HH11	1:A:312:ARG:HG3	1.52	0.75
2:B:338:LEU:N	2:B:338:LEU:HD23	2.01	0.75
1:A:437:THR:HG21	1:A:508:LEU:CD2	2.16	0.75
1:A:504:LEU:N	1:A:504:LEU:HD23	2.00	0.75
1:A:793:ILE:CG2	1:A:828:GLN:CD	2.60	0.74
1:A:511:LEU:N	1:A:511:LEU:HD23	2.01	0.74
1:A:494:TYR:CD2	2:B:367:ILE:HD13	2.23	0.74
1:A:198:ASP:OD1	1:A:198:ASP:N	2.17	0.74
1:A:793:ILE:HD12	1:A:793:ILE:N	2.02	0.74
1:A:407:SER:CB	1:A:545:SER:O	2.35	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:667:ASP:CB	1:A:744:LYS:NZ	2.51	0.73
2:B:425:ARG:HA	2:B:430:ILE:HG23	1.70	0.73
1:A:469:LYS:HA	1:A:469:LYS:CE	2.18	0.73
1:A:815:LEU:HD12	1:A:815:LEU:C	2.14	0.72
2:B:334:VAL:HG12	2:B:335:LEU:HD23	1.70	0.72
1:A:308:GLU:CD	3:A:901:YAO:O17	2.32	0.72
2:B:425:ARG:HG3	2:B:425:ARG:NH1	2.00	0.72
1:A:659:LEU:HD12	1:A:659:LEU:C	2.15	0.71
1:A:325:TYR:CE2	1:A:665:CYS:HB3	2.27	0.70
1:A:667:ASP:HB3	1:A:744:LYS:HZ3	1.53	0.70
1:A:700:ALA:HB1	1:A:701:PRO:HD2	1.72	0.69
1:A:538:PHE:CE1	1:A:706:LEU:HD21	2.26	0.69
1:A:750:ARG:HG3	1:A:750:ARG:NH1	2.02	0.69
2:B:395:ILE:HG22	2:B:433:VAL:HG11	1.75	0.69
1:A:320:PHE:CD1	1:A:747:VAL:HG21	2.27	0.69
2:B:337:GLN:N	2:B:337:GLN:OE1	2.25	0.69
2:B:399:GLY:N	2:B:437:TRP:CD1	2.61	0.69
1:A:308:GLU:OE2	3:A:901:YAO:O17	2.11	0.69
1:A:370:VAL:HG21	1:A:528:ILE:HD13	1.75	0.68
1:A:356:ILE:HD11	1:A:566:THR:HG23	1.76	0.68
1:A:793:ILE:HG23	1:A:828:GLN:CD	2.19	0.68
1:A:392:LEU:CD1	2:B:316:LEU:HD22	2.23	0.67
1:A:312:ARG:HG3	1:A:312:ARG:NH1	2.08	0.67
1:A:754:ASP:OD1	1:A:755:PRO:HD2	1.94	0.67
2:B:327:ASN:ND2	2:B:330:ALA:HB2	2.09	0.67
1:A:656:PHE:HE2	1:A:759:GLY:HA3	1.56	0.67
1:A:384:ARG:HH22	2:B:312:LYS:C	2.03	0.66
1:A:793:ILE:HG22	1:A:794:PRO:HD2	1.77	0.66
1:A:465:ALA:HB1	1:A:479:LEU:HD23	1.78	0.66
1:A:656:PHE:CE2	1:A:759:GLY:CA	2.76	0.66
1:A:437:THR:HG21	1:A:508:LEU:HD23	1.79	0.65
1:A:296:GLN:O	1:A:299:SER:HB3	1.96	0.65
2:B:395:ILE:HG22	2:B:433:VAL:CG1	2.26	0.64
1:A:349:VAL:HG21	1:A:577:ALA:HB1	1.78	0.64
2:B:398:TYR:C	2:B:437:TRP:NE1	2.55	0.64
2:B:363:LEU:HD23	2:B:363:LEU:N	2.13	0.64
2:B:399:GLY:H	2:B:437:TRP:CD1	2.16	0.64
1:A:695:TRP:NE1	1:A:706:LEU:CD1	2.61	0.63
1:A:484:HIS:CE1	2:B:372:LEU:HD13	2.34	0.63
2:B:398:TYR:C	2:B:437:TRP:HE1	2.05	0.63
1:A:658:ASN:HA	1:A:760:SER:CB	2.30	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:486:ASP:OD1	2:B:398:TYR:OH	2.17	0.62
1:A:407:SER:HB2	1:A:545:SER:O	2.00	0.62
1:A:352:GLU:HB3	1:A:568:ARG:HB2	1.82	0.62
1:A:485:ARG:HE	2:B:404:ALA:CB	2.13	0.61
1:A:510:GLU:HG2	1:A:511:LEU:HD23	1.82	0.61
1:A:726:ARG:O	1:A:730:ILE:HG13	1.99	0.61
1:A:656:PHE:CD2	1:A:759:GLY:HA3	2.35	0.61
1:A:320:PHE:CE1	1:A:747:VAL:CG2	2.83	0.61
1:A:530:ASP:OD2	1:A:685:THR:HG22	2.01	0.60
1:A:752:ARG:HG3	1:A:752:ARG:HH11	1.66	0.60
1:A:568:ARG:NH2	1:A:699:LYS:HG2	2.17	0.60
1:A:804:ILE:HG23	1:A:804:ILE:O	2.01	0.60
1:A:370:VAL:HG21	1:A:528:ILE:CD1	2.32	0.60
1:A:317:VAL:HG12	1:A:317:VAL:O	2.02	0.59
1:A:667:ASP:CB	1:A:744:LYS:HZ3	2.13	0.59
1:A:667:ASP:OD2	1:A:744:LYS:NZ	2.29	0.59
1:A:349:VAL:HG12	1:A:349:VAL:O	2.03	0.59
2:B:399:GLY:HA3	2:B:437:TRP:CE2	2.37	0.59
1:A:801:GLU:HG3	1:A:809:ALA:CA	2.27	0.59
2:B:370:TYR:HD2	2:B:370:TYR:N	2.01	0.58
1:A:349:VAL:HG21	1:A:577:ALA:CB	2.33	0.58
1:A:661:LYS:HG3	1:A:704:LEU:HD23	1.86	0.58
1:A:695:TRP:CE3	1:A:697:LEU:HD11	2.38	0.58
1:A:326:VAL:HG12	1:A:326:VAL:O	2.03	0.58
1:A:537:GLU:HG2	1:A:544:LEU:CD2	2.34	0.57
1:A:538:PHE:CD1	1:A:706:LEU:CD2	2.87	0.57
1:A:538:PHE:CD1	1:A:706:LEU:HD23	2.39	0.57
2:B:337:GLN:O	2:B:341:GLU:N	2.32	0.57
1:A:192:GLU:CD	1:A:214:ARG:NH1	2.62	0.57
2:B:370:TYR:N	2:B:370:TYR:CD2	2.72	0.57
1:A:538:PHE:CE1	1:A:706:LEU:CD2	2.87	0.57
1:A:776:MET:HE2	1:A:802:HIS:HD2	1.70	0.57
2:B:432:GLU:O	2:B:435:GLN:HG2	2.04	0.57
1:A:353:LEU:HB3	1:A:565:LEU:HD23	1.86	0.57
1:A:793:ILE:H	1:A:793:ILE:CD1	2.02	0.57
1:A:801:GLU:HG2	1:A:809:ALA:H	1.70	0.57
1:A:363:TYR:CD2	1:A:734:ILE:HG13	2.40	0.57
2:B:317:SER:O	2:B:321:VAL:HG23	2.05	0.56
1:A:188:MET:HE3	1:A:210:PHE:CE2	2.40	0.56
1:A:533:PHE:O	1:A:537:GLU:HG3	2.05	0.56
1:A:645:GLU:O	1:A:645:GLU:HG3	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:667:ASP:CB	1:A:744:LYS:HZ1	2.15	0.56
1:A:574:VAL:HB	1:A:575:PRO:HD3	1.88	0.56
1:A:658:ASN:HA	1:A:760:SER:HB3	1.86	0.56
2:B:398:TYR:CA	2:B:437:TRP:HE1	2.19	0.56
1:A:684:THR:HG22	1:A:686:ALA:H	1.70	0.56
1:A:801:GLU:HG2	1:A:801:GLU:O	2.04	0.56
1:A:442:LYS:HE3	2:B:355:THR:HG21	1.88	0.56
1:A:752:ARG:HG3	1:A:752:ARG:NH1	2.19	0.56
1:A:793:ILE:HG22	1:A:794:PRO:CD	2.36	0.55
2:B:430:ILE:O	2:B:434:LEU:HD12	2.05	0.55
1:A:193:ALA:C	1:A:200:ILE:HD11	2.29	0.55
2:B:420:PHE:CD2	2:B:420:PHE:C	2.84	0.55
1:A:458:LEU:HB3	1:A:487:LEU:HD12	1.88	0.55
1:A:320:PHE:C	1:A:320:PHE:CD2	2.85	0.55
2:B:383:TRP:CZ2	2:B:420:PHE:HD1	2.25	0.55
1:A:501:GLN:O	1:A:505:GLU:HG3	2.08	0.54
1:A:441:LEU:HD23	2:B:356:ASN:HD22	1.72	0.54
1:A:465:ALA:CB	1:A:479:LEU:HD23	2.37	0.54
2:B:369:PRO:HB2	2:B:370:TYR:HD2	1.72	0.54
2:B:419:ASN:H	2:B:419:ASN:ND2	2.06	0.53
1:A:484:HIS:ND1	2:B:372:LEU:CD1	2.72	0.53
1:A:671:TRP:O	1:A:673:PRO:HD3	2.08	0.53
1:A:438:GLN:OE1	1:A:508:LEU:HD11	2.01	0.53
2:B:403:GLN:OE1	2:B:403:GLN:HA	2.06	0.53
1:A:718:ILE:HG22	1:A:723:ILE:HG13	1.91	0.53
2:B:419:ASN:N	2:B:419:ASN:HD22	2.05	0.53
1:A:332:MET:HG2	1:A:333:VAL:HG23	1.91	0.53
2:B:369:PRO:HB2	2:B:370:TYR:CD2	2.44	0.52
1:A:568:ARG:HH12	1:A:699:LYS:HG2	1.70	0.52
1:A:205:GLN:O	1:A:209:VAL:HG23	2.10	0.52
1:A:297:LEU:HB2	1:A:304:VAL:HG21	1.91	0.52
1:A:437:THR:CB	1:A:508:LEU:HD21	2.39	0.52
1:A:456:LYS:CB	2:B:370:TYR:CE1	2.84	0.52
1:A:740:VAL:O	1:A:740:VAL:HG13	2.09	0.51
2:B:368:GLU:N	2:B:369:PRO:CD	2.73	0.51
2:B:363:LEU:N	2:B:363:LEU:CD2	2.72	0.51
2:B:374:GLU:O	2:B:376:ILE:HD12	2.09	0.51
2:B:419:ASN:ND2	2:B:419:ASN:N	2.59	0.51
1:A:484:HIS:CE1	2:B:372:LEU:CD1	2.93	0.51
1:A:656:PHE:HE2	1:A:759:GLY:CA	2.18	0.51
1:A:658:ASN:HA	1:A:760:SER:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:TYR:CD2	1:A:734:ILE:HG23	2.46	0.51
1:A:485:ARG:HE	2:B:404:ALA:HB1	1.75	0.51
1:A:319:THR:HB	1:A:572:SER:CB	2.26	0.50
1:A:568:ARG:CZ	1:A:699:LYS:CG	2.87	0.50
3:A:901:YAO:O1	3:A:901:YAO:O3	2.28	0.50
1:A:724:VAL:O	1:A:728:LEU:HG	2.12	0.50
1:A:231:PHE:CE1	1:A:249:VAL:CG1	2.72	0.50
1:A:340:ASN:HB2	1:A:560:PHE:CD2	2.47	0.50
1:A:363:TYR:HB3	1:A:734:ILE:HG13	1.93	0.50
1:A:695:TRP:NE1	1:A:706:LEU:HD13	2.17	0.50
1:A:257:GLU:HG3	1:A:263:ASN:HD22	1.77	0.49
1:A:363:TYR:CE2	1:A:734:ILE:HG23	2.47	0.49
2:B:361:GLU:O	2:B:361:GLU:HG3	2.12	0.49
1:A:432:LYS:HA	1:A:435:VAL:HG22	1.94	0.49
1:A:695:TRP:HB3	1:A:697:LEU:HG	1.94	0.49
1:A:601:GLU:HA	1:A:616:TYR:O	2.13	0.49
1:A:392:LEU:HD23	1:A:398:PHE:CD2	2.48	0.48
1:A:391:TYR:CZ	2:B:310:PRO:HG3	2.49	0.48
1:A:422:HIS:CD2	2:B:315:PHE:CD2	3.02	0.48
1:A:271:LYS:HG3	1:A:271:LYS:O	2.13	0.48
1:A:458:LEU:HB3	1:A:487:LEU:CD1	2.44	0.48
1:A:261:LEU:HD23	1:A:261:LEU:HA	1.71	0.48
1:A:321:ARG:NH2	1:A:569:ASN:O	2.47	0.48
1:A:370:VAL:HG22	1:A:528:ILE:HD11	1.96	0.48
1:A:442:LYS:N	2:B:356:ASN:HD21	2.11	0.48
1:A:594:ARG:HA	1:A:640:VAL:O	2.14	0.47
1:A:662:VAL:HG13	1:A:748:VAL:HG22	1.96	0.47
1:A:694:PHE:HA	1:A:704:LEU:O	2.14	0.47
1:A:707:VAL:HG12	1:A:712:ALA:HA	1.95	0.47
1:A:422:HIS:NE2	2:B:315:PHE:CE2	2.82	0.47
1:A:437:THR:HG22	1:A:508:LEU:CD2	2.24	0.47
1:A:337:LEU:HD23	1:A:337:LEU:N	2.30	0.47
1:A:506:GLU:O	1:A:509:GLN:N	2.46	0.47
1:A:694:PHE:CE2	1:A:731:LEU:HD11	2.50	0.47
1:A:363:TYR:CB	1:A:734:ILE:HG13	2.44	0.47
1:A:370:VAL:CG2	1:A:528:ILE:HD11	2.44	0.47
1:A:312:ARG:NH1	1:A:312:ARG:CG	2.72	0.47
1:A:458:LEU:HD23	1:A:458:LEU:HA	1.69	0.46
2:B:359:LEU:HD23	2:B:359:LEU:H	1.79	0.46
1:A:238:LEU:HB2	1:A:243:ASN:HB3	1.96	0.46
2:B:399:GLY:CA	2:B:437:TRP:CE2	2.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:444:LEU:HD22	1:A:500:THR:HG22	1.98	0.46
1:A:663:VAL:CG2	1:A:747:VAL:HB	2.46	0.46
1:A:752:ARG:HH11	1:A:752:ARG:CG	2.29	0.46
1:A:370:VAL:CG2	1:A:528:ILE:CD1	2.93	0.46
1:A:780:ILE:HB	1:A:796:LEU:HB3	1.96	0.46
1:A:465:ALA:CB	1:A:479:LEU:CD2	2.93	0.46
1:A:548:SER:O	1:A:552:TRP:HB3	2.15	0.46
1:A:188:MET:HG2	1:A:210:PHE:HE2	1.80	0.46
1:A:542:THR:OG1	1:A:543:PRO:HD2	2.16	0.46
1:A:238:LEU:CB	1:A:243:ASN:HB3	2.46	0.46
1:A:583:ASP:OD1	1:A:585:LYS:NZ	2.43	0.45
1:A:235:LEU:HD21	1:A:246:THR:HG22	1.97	0.45
1:A:350:ASN:ND2	1:A:350:ASN:C	2.72	0.45
1:A:647:LYS:HE3	1:A:798:PHE:CE1	2.51	0.45
1:A:701:PRO:HG2	1:A:701:PRO:O	2.16	0.45
2:B:380:ASN:CG	2:B:381:ALA:H	2.24	0.45
1:A:680:HIS:CE1	1:A:730:ILE:HD13	2.51	0.45
2:B:416:GLN:HA	2:B:419:ASN:HD21	1.82	0.45
1:A:465:ALA:HB1	1:A:479:LEU:CD2	2.47	0.45
1:A:469:LYS:CE	1:A:469:LYS:CA	2.92	0.45
1:A:294:ALA:HA	1:A:304:VAL:HG11	1.99	0.45
1:A:521:LEU:HD22	1:A:525:ASP:HB3	1.99	0.45
2:B:350:GLN:O	2:B:354:GLN:HB2	2.17	0.44
1:A:474:ILE:HG23	2:B:393:GLN:NE2	2.31	0.44
1:A:720:ASP:O	1:A:724:VAL:HG23	2.17	0.44
1:A:209:VAL:HG12	1:A:213:ILE:HD11	1.99	0.44
1:A:188:MET:SD	1:A:200:ILE:HG23	2.57	0.44
1:A:321:ARG:HH11	1:A:321:ARG:HD3	1.66	0.44
1:A:378:VAL:O	1:A:381:GLU:N	2.51	0.44
1:A:661:LYS:HZ3	1:A:704:LEU:HD21	1.82	0.44
1:A:361:PRO:HG2	1:A:678:PHE:HB2	1.99	0.44
1:A:684:THR:HG22	1:A:685:THR:N	2.33	0.44
1:A:776:MET:HE2	1:A:802:HIS:CD2	2.52	0.44
1:A:789:ALA:HB1	1:A:790:PRO:HD2	2.00	0.44
1:A:188:MET:HG2	1:A:210:PHE:CE2	2.53	0.43
1:A:229:LEU:HA	1:A:229:LEU:HD12	1.85	0.43
2:B:337:GLN:OE1	2:B:337:GLN:CA	2.65	0.43
1:A:456:LYS:HG3	2:B:370:TYR:CE1	2.53	0.43
1:A:657:GLY:HA2	1:A:752:ARG:NH2	2.33	0.43
1:A:793:ILE:CG2	1:A:794:PRO:CD	2.96	0.43
1:A:644:PRO:HD2	1:A:780:ILE:CD1	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:695:TRP:NE1	1:A:706:LEU:HD11	2.31	0.43
1:A:791:GLN:HE21	1:A:791:GLN:HB3	1.67	0.43
1:A:815:LEU:C	1:A:815:LEU:CD1	2.84	0.43
2:B:367:ILE:O	2:B:367:ILE:HG13	2.18	0.43
1:A:658:ASN:ND2	1:A:752:ARG:HB2	2.34	0.43
2:B:419:ASN:H	2:B:419:ASN:HD22	1.66	0.43
1:A:492:LYS:HD2	1:A:492:LYS:HA	1.67	0.43
1:A:283:ILE:HD12	1:A:294:ALA:HB2	2.00	0.43
1:A:221:TRP:CD1	1:A:262:ILE:HA	2.54	0.43
1:A:754:ASP:OD1	1:A:755:PRO:CD	2.65	0.43
1:A:503:LYS:HB2	1:A:504:LEU:HD23	2.01	0.42
1:A:222:LEU:HD23	1:A:222:LEU:HA	1.87	0.42
1:A:469:LYS:HE2	1:A:470:PRO:HD3	2.01	0.42
1:A:463:LYS:HD2	1:A:463:LYS:HA	1.90	0.42
1:A:734:ILE:HD12	1:A:734:ILE:N	2.34	0.42
2:B:425:ARG:HH11	2:B:425:ARG:CG	2.11	0.42
1:A:484:HIS:ND1	2:B:372:LEU:HD13	2.34	0.42
1:A:456:LYS:CG	2:B:370:TYR:HE1	2.31	0.42
1:A:456:LYS:HB2	2:B:370:TYR:HE1	1.82	0.42
1:A:537:GLU:HG2	1:A:544:LEU:HD21	2.01	0.42
1:A:297:LEU:HA	1:A:297:LEU:HD23	1.75	0.42
1:A:192:GLU:CD	1:A:214:ARG:HH12	2.28	0.42
1:A:235:LEU:HD13	1:A:249:VAL:HG21	2.01	0.42
1:A:494:TYR:OH	2:B:363:LEU:CD1	2.68	0.42
1:A:798:PHE:O	1:A:821:GLU:HG3	2.20	0.42
1:A:830:LEU:N	1:A:830:LEU:HD23	2.33	0.42
1:A:374:LYS:HZ3	1:A:374:LYS:HG3	1.71	0.41
1:A:439:GLU:OE2	2:B:352:ILE:HD11	2.20	0.41
2:B:324:VAL:HG13	2:B:331:ALA:CB	2.50	0.41
1:A:335:THR:O	1:A:335:THR:OG1	2.23	0.41
1:A:654:MET:HE2	1:A:654:MET:HB3	1.89	0.41
1:A:180:GLN:HA	1:A:339:GLY:HA2	2.01	0.41
1:A:422:HIS:NE2	2:B:315:PHE:CZ	2.89	0.41
1:A:451:LEU:HD23	1:A:494:TYR:HB2	2.02	0.41
1:A:793:ILE:CG2	1:A:794:PRO:HD2	2.49	0.41
2:B:388:GLN:O	2:B:391:ALA:HB3	2.20	0.41
1:A:253:HIS:ND1	1:A:253:HIS:C	2.78	0.41
2:B:383:TRP:CE3	2:B:412:LYS:HE2	2.55	0.41
2:B:377:GLN:OE1	2:B:411:ASN:HB3	2.20	0.41
2:B:402:PHE:HB2	2:B:414:VAL:HG13	2.02	0.41
2:B:402:PHE:CE2	2:B:418:LYS:HG3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:TYR:CG	1:A:734:ILE:HG13	2.55	0.41
2:B:383:TRP:CE2	2:B:420:PHE:HD1	2.39	0.41
2:B:425:ARG:HA	2:B:430:ILE:CG2	2.44	0.41
1:A:311:ASP:OD1	1:A:311:ASP:N	2.43	0.40
1:A:485:ARG:NE	2:B:404:ALA:HB1	2.36	0.40
1:A:601:GLU:HB3	1:A:617:LYS:HD3	2.02	0.40
1:A:657:GLY:C	1:A:752:ARG:NH2	2.80	0.40
1:A:659:LEU:O	1:A:659:LEU:HG	2.21	0.40
1:A:715:MET:HE3	1:A:723:ILE:HG12	2.02	0.40
1:A:791:GLN:HA	1:A:792:PRO:HD3	1.82	0.40
1:A:521:LEU:HD23	1:A:521:LEU:HA	1.94	0.40
1:A:353:LEU:HB3	1:A:565:LEU:CD2	2.51	0.40
1:A:665:CYS:O	1:A:744:LYS:N	2.51	0.40
1:A:297:LEU:CB	1:A:304:VAL:HG21	2.52	0.40
1:A:353:LEU:HA	1:A:353:LEU:HD23	1.67	0.40
1:A:370:VAL:HG12	1:A:375:ASP:HB2	2.02	0.40
1:A:603:ILE:HG12	1:A:615:ILE:CD1	2.51	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	664/871 (76%)	639 (96%)	24 (4%)	1 (0%)	43	70
2	B	131/144 (91%)	121 (92%)	10 (8%)	0	100	100
All	All	795/1015 (78%)	760 (96%)	34 (4%)	1 (0%)	48	76

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	338	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	566/715 (79%)	524 (93%)	42 (7%)	13	36
2	B	117/125 (94%)	94 (80%)	23 (20%)	1	6
All	All	683/840 (81%)	618 (90%)	65 (10%)	8	27

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	191	GLN
1	A	198	ASP
1	A	200	ILE
1	A	249	VAL
1	A	319	THR
1	A	320	PHE
1	A	326	VAL
1	A	349	VAL
1	A	350	ASN
1	A	352	GLU
1	A	353	LEU
1	A	370	VAL
1	A	373	GLU
1	A	374	LYS
1	A	380	GLN
1	A	402	ASN
1	A	404	LYS
1	A	458	LEU
1	A	471	PRO
1	A	472	ARG
1	A	488	THR
1	A	490	LEU
1	A	503	LYS
1	A	504	LEU
1	A	505	GLU
1	A	508	LEU
1	A	509	GLN

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Mol	Chain	Res	Type
1	A	511	LEU
1	A	569	ASN
1	A	659	LEU
1	A	661	LYS
1	A	663	VAL
1	A	704	LEU
1	A	731	LEU
1	A	780	ILE
1	A	786	ILE
1	A	791	GLN
1	A	793	ILE
1	A	801	GLU
1	A	811	VAL
1	A	815	LEU
1	A	824	ARG
2	B	332	THR
2	B	334	VAL
2	B	335	LEU
2	B	336	ARG
2	B	337	GLN
2	B	338	LEU
2	B	359	LEU
2	B	363	LEU
2	B	367	ILE
2	B	368	GLU
2	B	370	TYR
2	B	371	ARG
2	B	372	LEU
2	B	405	ILE
2	B	415	VAL
2	B	419	ASN
2	B	420	PHE
2	B	422	VAL
2	B	425	ARG
2	B	427	ARG
2	B	430	ILE
2	B	436	GLU
2	B	438	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	350	ASN
1	A	380	GLN
1	A	402	ASN
1	A	509	GLN
1	A	514	ASN
1	A	551	HIS
1	A	633	GLN
1	A	778	GLN
1	A	791	GLN
2	B	327	ASN
2	B	356	ASN
2	B	419	ASN
2	B	423	ASN

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 ⓘ

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	YAO	A	901	-	77,80,80	1.96	20 (25%)	105,123,123	1.42	14 (13%)

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	YAO	A	901	-	-	8/46/114/114	0/8/9/9

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	901	YAO	P1-O11	-5.28	1.53	1.59
3	A	901	YAO	P2-O11	-5.14	1.53	1.59
3	A	901	YAO	C5-N1	-4.72	1.32	1.41
3	A	901	YAO	O5-C30	-4.70	1.33	1.43
3	A	901	YAO	C37-N6	-4.58	1.29	1.37
3	A	901	YAO	C26-N3	-4.18	1.30	1.37
3	A	901	YAO	C27-N4	-3.49	1.28	1.36
3	A	901	YAO	C36-N6	-3.33	1.37	1.46
3	A	901	YAO	O6-C31	-3.32	1.34	1.43
3	A	901	YAO	C9-N1	-3.22	1.41	1.46
3	A	901	YAO	C5-C4	-2.92	1.35	1.41
3	A	901	YAO	C6-C7	-2.69	1.36	1.39
3	A	901	YAO	C29-C30	-2.60	1.49	1.52
3	A	901	YAO	C28-N5	-2.58	1.32	1.38
3	A	901	YAO	O1-C9	-2.45	1.36	1.41
3	A	901	YAO	P2-O12	-2.45	1.44	1.55
3	A	901	YAO	O4-C27	-2.41	1.19	1.24
3	A	901	YAO	P1-O9	-2.35	1.44	1.55
3	A	901	YAO	C3-C2	-2.34	1.36	1.39
3	A	901	YAO	C27-N3	-2.08	1.34	1.39

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	YAO	C25-C11-C12	5.61	124.79	116.08
3	A	901	YAO	C5-N1-C9	-5.00	112.46	123.92
3	A	901	YAO	C11-C10-C9	-3.89	96.41	104.25
3	A	901	YAO	C26-C25-C28	-3.87	106.53	113.40
3	A	901	YAO	C26-N3-C27	-3.50	120.14	125.42
3	A	901	YAO	O6-C31-C32	-3.15	101.76	108.93
3	A	901	YAO	O15-C36-C42	-2.93	100.35	106.62
3	A	901	YAO	C43-C42-C36	-2.87	96.04	101.46
3	A	901	YAO	O4-C27-N4	-2.23	118.09	121.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	YAO	O17-C43-C42	-2.15	104.93	111.82
3	A	901	YAO	O7-C32-C31	-2.11	104.32	109.25
3	A	901	YAO	N3-C27-N4	2.04	123.84	119.50
3	A	901	YAO	O11-P1-O10	-2.03	104.61	110.70
3	A	901	YAO	O8-P1-O10	-2.01	100.99	108.94

There are no chirality outliers.

All (8) torsion outliers are listed below:

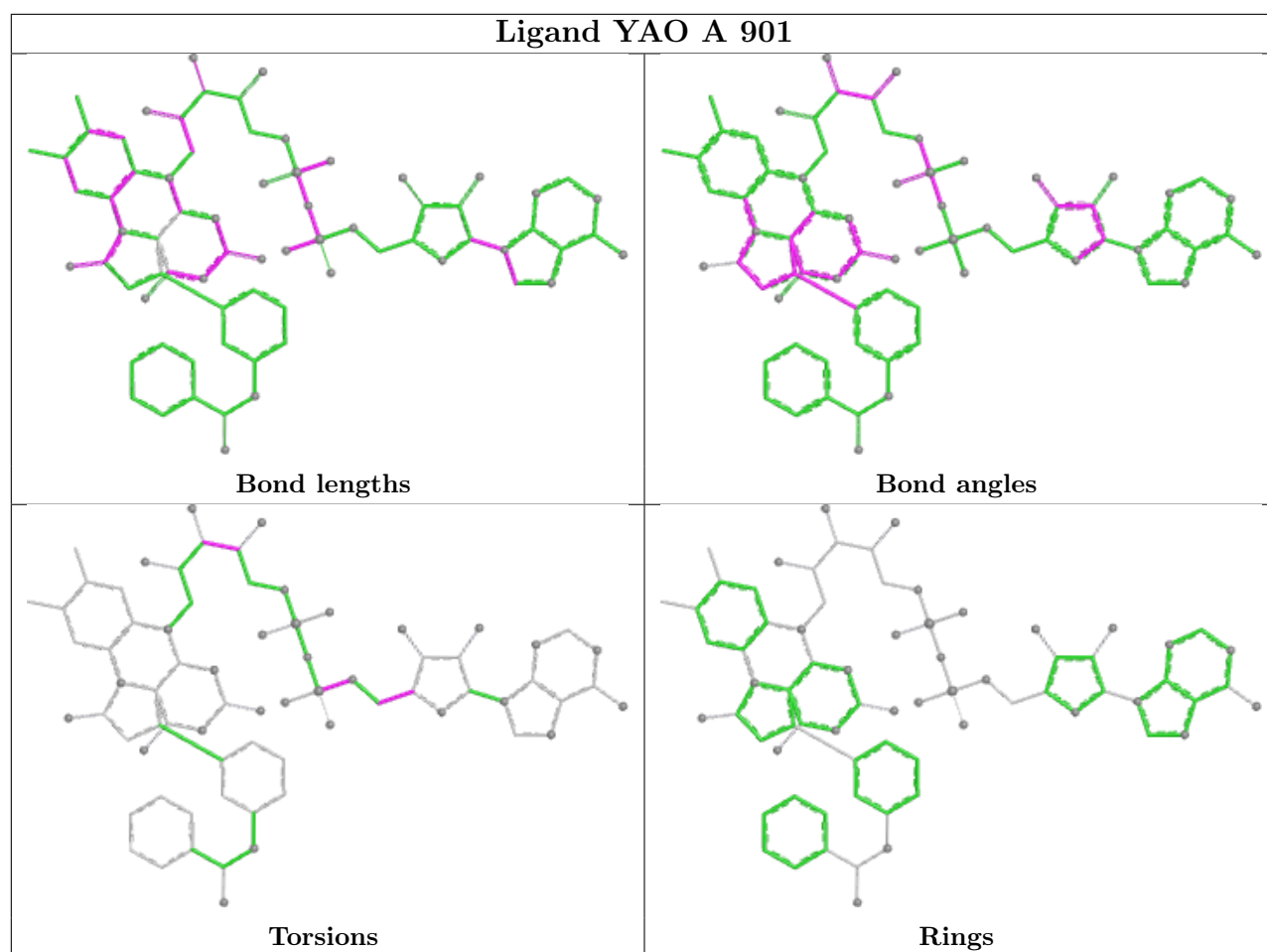
Mol	Chain	Res	Type	Atoms
3	A	901	YAO	C34-O14-P2-O11
3	A	901	YAO	C34-O14-P2-O13
3	A	901	YAO	O14-C34-C35-C43
3	A	901	YAO	O6-C31-C32-C33
3	A	901	YAO	O14-C34-C35-O15
3	A	901	YAO	O6-C31-C32-O7
3	A	901	YAO	C34-O14-P2-O12
3	A	901	YAO	C30-C31-C32-O7

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	YAO	4	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 [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	666/871 (76%)	0.02	5 (0%) 82 65	67, 104, 138, 162	0
2	B	133/144 (92%)	0.31	2 (1%) 72 51	97, 134, 155, 167	0
All	All	799/1015 (78%)	0.07	7 (0%) 81 63	67, 109, 146, 167	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	836	LEU	5.1
1	A	171	PRO	5.0
1	A	377	MET	2.5
2	B	337	GLN	2.2
1	A	785	SER	2.1
1	A	683	SER	2.0
2	B	427	ARG	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](#)

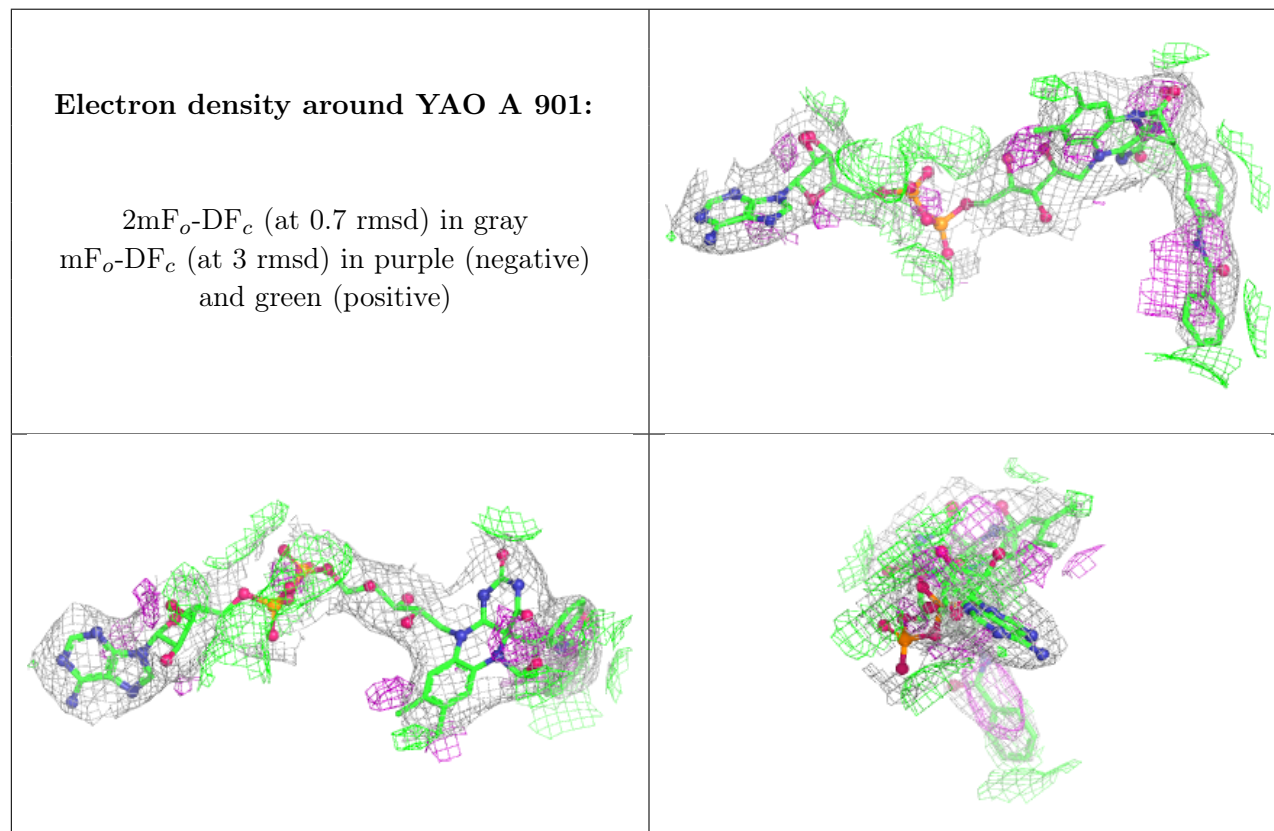
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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
3	YAO	A	901	72/72	0.94	0.12	68,86,114,124	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.