



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 08:25 PM UTC

PDB ID : 8F59 / pdb_00008f59
Title : LSD1-CoREST in complex with AW2 and SNAG peptide
Authors : Caroli, J.; Mattevi, A.
Deposited on : 2022-11-12
Resolution : 2.80 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

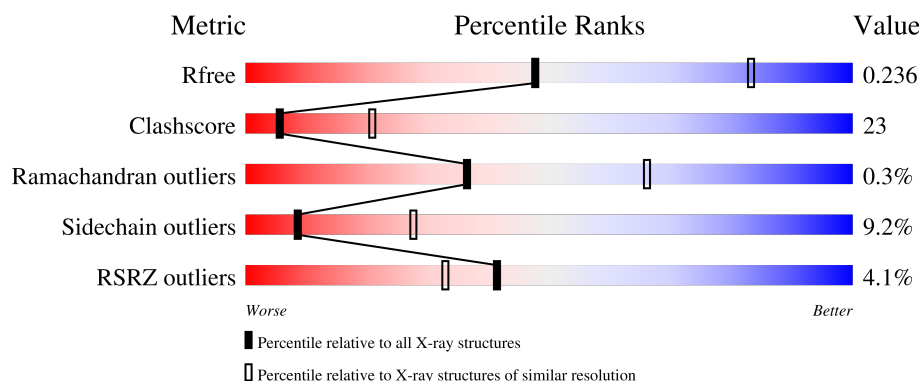
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	2% 54% 19% • 24%
2	B	144	6% 48% 30% 13% • 8%
3	C	9	44% 44% 56%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6397 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			

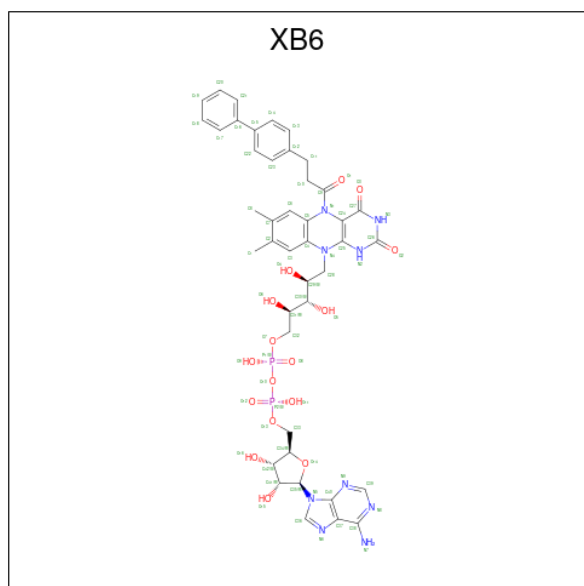
There are 8 discrepancies between the modelled and reference sequences:

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 a protein called Zinc finger protein SNAI1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	4	Total	C	N	O	0	0	0
			35	23	7	5			

- Molecule 4 is [(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-3,4-dihydroxyoxolan-2-yl]methyl (2R,3S,4S)-5-{5-[3-([1,1'-biphenyl]-4-yl)propanoyl]-7,8-dimethyl-2,4-dioxo-1,3,4,5-tetrahydrobenzo[g]pteridin-10(2H)-yl}-2,3,4-trihydroxypentyl dihydrogen diphosphate (non-preferred name) (CCD ID: XB6) (formula: C₄₂H₄₇N₉O₁₆P₂) (labeled as "Ligand of Interest" by depositor).

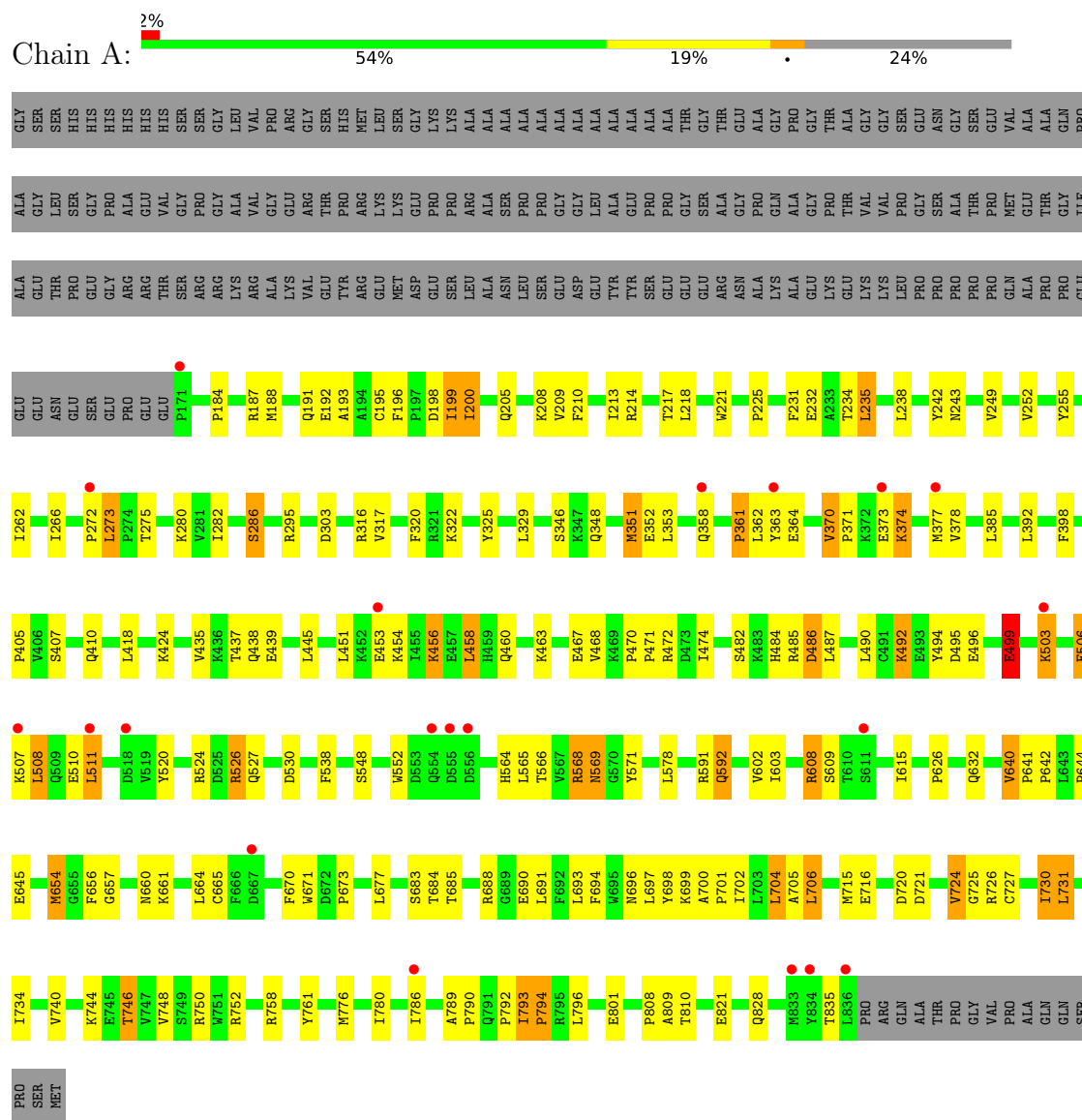


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			69	42	9	16	2		

3 Residue-property plots

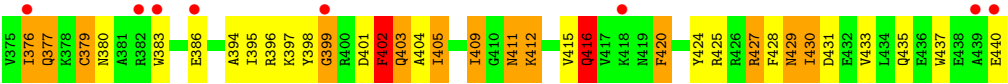
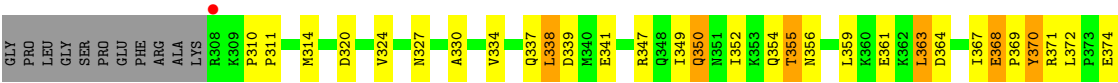
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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

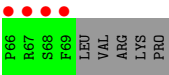


• Molecule 2: REST corepressor 1





● Molecule 3: Zinc finger protein SNAI1



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	120.21Å 179.20Å 236.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.73 – 2.80 48.73 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.73-2.80) 99.4 (48.73-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.18.2	Depositor
R, R_{free}	0.221 , 0.234 0.227 , 0.236	Depositor DCC
R_{free} test set	1233 reflections (1.96%)	wwPDB-VP
Wilson B-factor (Å ²)	90.9	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 70.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6397	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.16% 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

Bond lengths and bond angles in the following residue types are not validated in this section: XB6

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	0.85	6/5331 (0.1%)	1.09	9/7232 (0.1%)
2	B	0.90	1/1091 (0.1%)	1.42	2/1471 (0.1%)
3	C	0.67	0/36	0.58	0/46
All	All	0.85	7/6458 (0.1%)	1.15	11/8749 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	405	PRO	C-O	-6.00	1.17	1.23
1	A	644	PRO	C-O	-5.86	1.16	1.23
1	A	184	PRO	C-O	-5.65	1.17	1.23
1	A	361	PRO	C-O	-5.41	1.17	1.24
1	A	706	LEU	C-O	-5.31	1.17	1.23
2	B	310	PRO	CA-C	-5.21	1.49	1.51
1	A	792	PRO	C-O	-5.05	1.17	1.24

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	416	GLN	CB-CA-C	6.86	123.85	110.67
1	A	746	THR	CB-CA-C	-5.79	99.22	109.70
1	A	499	GLU	N-CA-C	-5.74	104.66	111.03
1	A	507	LYS	N-CA-C	-5.56	105.12	111.07
2	B	411	ASN	N-CA-C	-5.45	107.89	114.75
1	A	378	VAL	CA-C-O	-5.18	115.74	121.29
1	A	272	PRO	N-CA-C	-5.14	102.62	110.95
1	A	642	PRO	CB-CA-C	-5.11	104.69	111.23
1	A	794	PRO	N-CD-CG	-5.08	95.58	103.20
1	A	654	MET	CA-C-N	-5.05	117.90	122.80
1	A	654	MET	C-N-CA	-5.05	117.90	122.80

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	226	0
2	B	1076	0	1091	95	0
3	C	35	0	34	0	0
4	A	69	0	0	3	0
All	All	6397	0	6377	300	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 23.

All (300) 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:492:LYS:HA	1:A:492:LYS:HE2	1.21	1.13
1:A:363:TYR:HD2	1:A:734:ILE:HG13	1.12	1.07
1:A:438:GLN:OE1	1:A:508:LEU:HD12	1.54	1.07
1:A:438:GLN:OE1	1:A:508:LEU:CD1	2.00	1.07
1:A:608:ARG:HG3	1:A:608:ARG:HH11	1.18	1.04
1:A:363:TYR:CE2	1:A:734:ILE:HG23	1.94	1.02
1:A:458:LEU:HD12	1:A:487:LEU:HD12	1.41	1.02
1:A:346:SER:HA	1:A:351:MET:HE2	1.44	0.99
1:A:526:ARG:HH21	1:A:526:ARG:HG3	1.27	0.98
1:A:188:MET:CE	1:A:200:ILE:HB	1.94	0.98
1:A:485:ARG:NH1	2:B:404:ALA:HB1	1.78	0.97
1:A:363:TYR:CD2	1:A:734:ILE:HG13	1.99	0.95
1:A:720:ASP:O	1:A:724:VAL:HG23	1.67	0.95
1:A:363:TYR:HD2	1:A:734:ILE:CG1	1.79	0.94
1:A:458:LEU:CD1	1:A:487:LEU:HD12	1.98	0.93
2:B:376:ILE:HD12	2:B:376:ILE:H	1.32	0.93
1:A:492:LYS:HA	1:A:492:LYS:CE	1.99	0.91
2:B:311:PRO:HG2	2:B:314:MET:HG3	1.50	0.90
2:B:383:TRP:CZ2	2:B:416:GLN:HG2	2.06	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:346:SER:HA	1:A:351:MET:CE	2.01	0.89
1:A:654:MET:HE1	1:A:776:MET:HG3	1.51	0.89
1:A:188:MET:CE	1:A:200:ILE:CB	2.49	0.89
1:A:188:MET:CE	1:A:200:ILE:HG13	2.02	0.88
1:A:188:MET:HE2	1:A:200:ILE:HG13	1.57	0.86
1:A:363:TYR:HE2	1:A:734:ILE:HG23	1.33	0.86
1:A:506:GLU:OE1	1:A:506:GLU:N	2.09	0.86
1:A:188:MET:HE1	1:A:200:ILE:HB	1.59	0.84
1:A:654:MET:HE1	1:A:776:MET:CG	2.08	0.84
1:A:438:GLN:OE1	1:A:508:LEU:HD11	1.78	0.82
1:A:235:LEU:O	1:A:235:LEU:HD12	1.79	0.82
2:B:402:PHE:H	2:B:402:PHE:HD2	1.25	0.81
1:A:569:ASN:OD1	1:A:569:ASN:N	2.12	0.80
1:A:793:ILE:HG22	1:A:794:PRO:CD	2.12	0.80
1:A:700:ALA:HB1	1:A:701:PRO:HD2	1.61	0.79
1:A:685:THR:O	1:A:688:ARG:HG2	1.84	0.78
1:A:690:GLU:OE1	1:A:726:ARG:NH1	2.17	0.78
1:A:445:LEU:HB2	2:B:359:LEU:HD23	1.64	0.78
2:B:431:ASP:O	2:B:435:GLN:HG3	1.85	0.77
1:A:568:ARG:NH1	1:A:699:LYS:N	2.34	0.76
1:A:363:TYR:CE2	1:A:734:ILE:CG2	2.68	0.76
1:A:435:VAL:HG12	2:B:349:ILE:HD13	1.68	0.76
2:B:383:TRP:CH2	2:B:416:GLN:HG2	2.19	0.76
1:A:418:LEU:HD21	2:B:324:VAL:HG21	1.67	0.76
2:B:425:ARG:HA	2:B:430:ILE:HD13	1.66	0.76
1:A:188:MET:CE	1:A:200:ILE:CG1	2.64	0.76
1:A:458:LEU:CD1	1:A:487:LEU:CD1	2.63	0.76
1:A:608:ARG:HG3	1:A:608:ARG:NH1	1.98	0.76
1:A:451:LEU:HD23	1:A:494:TYR:HB2	1.69	0.75
1:A:511:LEU:N	1:A:511:LEU:HD23	2.00	0.75
1:A:435:VAL:HG13	2:B:349:ILE:HD11	1.70	0.74
2:B:376:ILE:HD12	2:B:376:ILE:N	2.03	0.74
1:A:363:TYR:CD2	1:A:734:ILE:HG23	2.23	0.73
1:A:458:LEU:HD13	1:A:487:LEU:CD1	2.19	0.73
1:A:209:VAL:HG12	1:A:213:ILE:HD11	1.71	0.72
1:A:188:MET:HE3	1:A:200:ILE:HB	1.71	0.72
1:A:463:LYS:O	1:A:467:GLU:HG2	1.92	0.70
1:A:654:MET:CE	1:A:776:MET:CG	2.70	0.70
2:B:430:ILE:N	2:B:430:ILE:HD12	2.07	0.70
2:B:350:GLN:HA	2:B:350:GLN:HE21	1.56	0.69
2:B:395:ILE:HG22	2:B:433:VAL:CG1	2.20	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:458:LEU:HB3	1:A:487:LEU:CD1	2.21	0.69
1:A:209:VAL:O	1:A:213:ILE:HD12	1.93	0.69
1:A:435:VAL:CG1	2:B:349:ILE:CD1	2.70	0.69
2:B:394:ALA:O	2:B:398:TYR:N	2.26	0.69
1:A:526:ARG:HG3	1:A:526:ARG:NH2	2.06	0.69
1:A:458:LEU:HD13	1:A:487:LEU:HD13	1.75	0.68
2:B:338:LEU:N	2:B:338:LEU:HD23	2.07	0.68
1:A:435:VAL:HG12	2:B:349:ILE:CD1	2.23	0.68
1:A:346:SER:CA	1:A:351:MET:CE	2.72	0.68
1:A:437:THR:OG1	1:A:508:LEU:HD21	1.92	0.68
1:A:273:LEU:O	1:A:273:LEU:HD23	1.94	0.68
1:A:205:GLN:HA	1:A:205:GLN:NE2	2.09	0.68
1:A:486:ASP:OD1	1:A:486:ASP:N	2.24	0.68
1:A:485:ARG:HH12	2:B:404:ALA:HB1	1.58	0.67
1:A:661:LYS:HD3	1:A:704:LEU:HD21	1.75	0.67
1:A:485:ARG:HH22	2:B:398:TYR:HE2	1.43	0.67
2:B:376:ILE:H	2:B:376:ILE:CD1	2.03	0.67
2:B:409:ILE:HD12	2:B:409:ILE:N	2.11	0.66
2:B:403:GLN:NE2	2:B:403:GLN:O	2.28	0.66
1:A:363:TYR:CD2	1:A:734:ILE:CG1	2.69	0.66
1:A:188:MET:HE1	1:A:200:ILE:CB	2.22	0.65
2:B:395:ILE:CG2	2:B:433:VAL:HG11	2.26	0.65
1:A:225:PRO:O	1:A:348:GLN:NE2	2.29	0.65
1:A:568:ARG:NH1	1:A:698:TYR:C	2.55	0.64
1:A:205:GLN:HA	1:A:205:GLN:HE21	1.63	0.64
1:A:363:TYR:CD2	1:A:734:ILE:CG2	2.81	0.64
1:A:286:SER:O	1:A:286:SER:OG	2.15	0.63
1:A:793:ILE:HG22	1:A:794:PRO:HD2	1.80	0.63
1:A:392:LEU:HD21	1:A:398:PHE:CD2	2.34	0.63
2:B:350:GLN:HE21	2:B:350:GLN:CA	2.10	0.62
1:A:373:GLU:OE1	1:A:373:GLU:N	2.27	0.62
1:A:199:ILE:N	1:A:199:ILE:HD13	2.14	0.62
1:A:188:MET:HE3	1:A:200:ILE:HG13	1.77	0.62
1:A:526:ARG:HH21	1:A:526:ARG:CG	2.08	0.61
1:A:439:GLU:HG3	2:B:352:ILE:CD1	2.31	0.61
2:B:395:ILE:O	2:B:399:GLY:N	2.33	0.61
1:A:353:LEU:HB3	1:A:565:LEU:HD13	1.83	0.61
2:B:368:GLU:N	2:B:369:PRO:CD	2.63	0.61
2:B:380:ASN:HB3	2:B:411:ASN:OD1	2.01	0.61
1:A:740:VAL:HG12	1:A:740:VAL:O	2.01	0.61
1:A:499:GLU:HA	1:A:499:GLU:OE1	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:510:GLU:HG2	1:A:511:LEU:HD23	1.83	0.60
1:A:392:LEU:HD23	1:A:398:PHE:HB3	1.83	0.60
1:A:821:GLU:HA	1:A:821:GLU:OE1	2.02	0.60
2:B:337:GLN:N	2:B:337:GLN:OE1	2.35	0.59
1:A:456:LYS:HA	2:B:370:TYR:HE1	1.67	0.59
1:A:566:THR:HG21	1:A:697:LEU:HD22	1.83	0.59
2:B:367:ILE:O	2:B:367:ILE:HG13	2.03	0.59
2:B:395:ILE:CG2	2:B:433:VAL:CG1	2.80	0.58
1:A:235:LEU:HD11	1:A:243:ASN:HB2	1.84	0.58
1:A:320:PHE:HB2	1:A:329:LEU:HD11	1.84	0.58
1:A:362:LEU:N	1:A:362:LEU:HD12	2.19	0.58
1:A:188:MET:HE3	1:A:200:ILE:CG1	2.31	0.58
1:A:353:LEU:N	1:A:353:LEU:HD22	2.19	0.58
2:B:412:LYS:HD2	2:B:412:LYS:N	2.18	0.58
1:A:568:ARG:HH12	1:A:699:LYS:HA	1.69	0.57
1:A:282:ILE:HG21	1:A:602:VAL:HG21	1.86	0.57
1:A:503:LYS:O	1:A:503:LYS:HG3	2.05	0.57
1:A:188:MET:HE1	1:A:200:ILE:CA	2.35	0.56
2:B:370:TYR:N	2:B:370:TYR:CD2	2.71	0.56
2:B:370:TYR:N	2:B:370:TYR:HD2	2.03	0.56
1:A:266:ILE:HD11	1:A:578:LEU:HD23	1.87	0.56
2:B:409:ILE:HD12	2:B:409:ILE:H	1.70	0.56
1:A:188:MET:HE2	1:A:200:ILE:CG1	2.29	0.56
1:A:435:VAL:HG13	2:B:349:ILE:CD1	2.34	0.56
1:A:456:LYS:HA	2:B:370:TYR:CE1	2.40	0.56
1:A:456:LYS:HG2	2:B:370:TYR:HE1	1.69	0.56
2:B:327:ASN:OD1	2:B:330:ALA:N	2.39	0.56
2:B:430:ILE:HD12	2:B:430:ILE:H	1.70	0.56
1:A:453:GLU:N	1:A:453:GLU:OE1	2.39	0.56
1:A:445:LEU:CB	2:B:359:LEU:HD23	2.34	0.55
1:A:351:MET:HA	1:A:569:ASN:HD21	1.69	0.55
1:A:435:VAL:CG1	2:B:349:ILE:HD11	2.34	0.55
1:A:568:ARG:HH12	1:A:699:LYS:N	2.03	0.55
2:B:420:PHE:CD2	2:B:420:PHE:C	2.85	0.55
2:B:350:GLN:HA	2:B:350:GLN:NE2	2.20	0.55
1:A:232:GLU:CD	1:A:232:GLU:H	2.15	0.54
1:A:424:LYS:HE3	2:B:339:ASP:OD1	2.08	0.54
1:A:195:CYS:HA	1:A:835:THR:HG22	1.88	0.54
1:A:508:LEU:N	1:A:508:LEU:HD23	2.23	0.54
1:A:468:VAL:O	1:A:472:ARG:NH1	2.40	0.54
1:A:808:PRO:O	1:A:810:THR:HG23	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:405:ILE:O	2:B:409:ILE:HD13	2.07	0.54
1:A:720:ASP:O	1:A:724:VAL:CG2	2.51	0.53
1:A:568:ARG:HH12	1:A:699:LYS:CA	2.21	0.53
1:A:316:ARG:NH1	1:A:656:PHE:HZ	2.05	0.53
2:B:402:PHE:CD2	2:B:402:PHE:N	2.70	0.53
2:B:430:ILE:H	2:B:430:ILE:CD1	2.22	0.53
1:A:362:LEU:N	1:A:362:LEU:CD1	2.72	0.53
2:B:383:TRP:CH2	2:B:420:PHE:HB2	2.43	0.53
1:A:363:TYR:HE2	1:A:734:ILE:CG2	2.09	0.53
1:A:603:ILE:HG12	1:A:615:ILE:HD13	1.90	0.53
1:A:188:MET:CE	1:A:200:ILE:CA	2.87	0.52
1:A:346:SER:CB	1:A:351:MET:CE	2.87	0.52
1:A:373:GLU:H	1:A:373:GLU:CD	2.16	0.52
2:B:430:ILE:N	2:B:430:ILE:CD1	2.72	0.52
1:A:407:SER:OG	1:A:410:GLN:HG3	2.10	0.52
2:B:368:GLU:N	2:B:369:PRO:HD3	2.25	0.52
1:A:494:TYR:C	1:A:494:TYR:CD1	2.88	0.51
1:A:196:PHE:HB3	1:A:199:ILE:HG12	1.92	0.51
1:A:235:LEU:HD12	1:A:235:LEU:C	2.31	0.51
1:A:325:TYR:CE1	1:A:665:CYS:HB3	2.45	0.51
1:A:654:MET:CE	1:A:776:MET:HG2	2.39	0.51
2:B:428:PHE:HB2	2:B:430:ILE:HD11	1.92	0.51
1:A:761:TYR:CD2	1:A:809:ALA:HB1	2.46	0.51
1:A:780:ILE:HB	1:A:796:LEU:HB3	1.93	0.51
2:B:380:ASN:CB	2:B:411:ASN:OD1	2.60	0.50
2:B:395:ILE:HG21	2:B:433:VAL:HG11	1.91	0.50
2:B:383:TRP:CH2	2:B:416:GLN:CG	2.92	0.50
2:B:401:ASP:C	2:B:403:GLN:N	2.69	0.50
1:A:564:HIS:O	1:A:565:LEU:HD23	2.12	0.50
1:A:231:PHE:CD1	1:A:231:PHE:C	2.89	0.49
1:A:701:PRO:HG2	1:A:701:PRO:O	2.12	0.49
2:B:379:CYS:SG	2:B:380:ASN:N	2.85	0.49
1:A:238:LEU:HB3	1:A:243:ASN:HB3	1.95	0.49
1:A:793:ILE:HG22	1:A:794:PRO:HD3	1.91	0.49
1:A:508:LEU:N	1:A:508:LEU:CD2	2.75	0.49
1:A:609:SER:O	1:A:609:SER:OG	2.27	0.48
1:A:361:PRO:O	1:A:361:PRO:HG2	2.12	0.48
2:B:396:ARG:O	2:B:396:ARG:HD3	2.14	0.48
1:A:456:LYS:HG2	2:B:370:TYR:CE1	2.49	0.48
1:A:210:PHE:CE1	1:A:252:VAL:HG22	2.48	0.48
1:A:358:GLN:O	1:A:358:GLN:HG3	2.10	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:495:ASP:CG	2:B:371:ARG:HH12	2.22	0.48
2:B:361:GLU:C	2:B:363:LEU:H	2.21	0.48
2:B:420:PHE:C	2:B:420:PHE:HD2	2.20	0.48
1:A:661:LYS:HD3	1:A:704:LEU:CD2	2.40	0.47
2:B:403:GLN:NE2	2:B:403:GLN:C	2.72	0.47
1:A:460:GLN:NE2	1:A:460:GLN:HA	2.28	0.47
1:A:482:SER:O	1:A:486:ASP:OD1	2.31	0.47
1:A:231:PHE:CE1	1:A:249:VAL:HG12	2.49	0.47
1:A:295:ARG:HD3	1:A:578:LEU:O	2.13	0.47
1:A:438:GLN:HB3	2:B:352:ILE:HG21	1.97	0.47
1:A:210:PHE:HE1	1:A:252:VAL:HG22	1.78	0.47
1:A:664:LEU:O	1:A:702:ILE:HA	2.14	0.47
1:A:193:ALA:HB2	1:A:200:ILE:HD12	1.96	0.47
1:A:789:ALA:HB1	1:A:790:PRO:HD2	1.96	0.47
1:A:726:ARG:O	1:A:730:ILE:HG12	2.14	0.47
2:B:429:ASN:HD22	2:B:429:ASN:N	2.13	0.47
1:A:373:GLU:N	1:A:373:GLU:CD	2.72	0.47
1:A:217:THR:HG23	1:A:234:THR:CG2	2.45	0.46
1:A:209:VAL:HG13	1:A:242:TYR:HD1	1.81	0.46
1:A:316:ARG:NH1	1:A:656:PHE:CZ	2.84	0.46
1:A:654:MET:HE2	1:A:776:MET:HG2	1.98	0.46
1:A:188:MET:HE3	1:A:200:ILE:CB	2.34	0.46
1:A:485:ARG:NH2	2:B:398:TYR:CE2	2.83	0.46
2:B:397:LYS:HD3	2:B:398:TYR:CZ	2.51	0.46
1:A:188:MET:HE1	1:A:200:ILE:N	2.31	0.46
1:A:209:VAL:C	1:A:213:ILE:HD12	2.41	0.46
1:A:828:GLN:O	1:A:828:GLN:HG3	2.16	0.46
1:A:192:GLU:OE2	1:A:214:ARG:NH1	2.46	0.45
2:B:374:GLU:O	2:B:374:GLU:HG2	2.16	0.45
1:A:484:HIS:CD2	2:B:372:LEU:HD13	2.51	0.45
2:B:377:GLN:O	2:B:377:GLN:HG3	2.09	0.45
2:B:395:ILE:HG22	2:B:433:VAL:HG11	1.88	0.45
1:A:608:ARG:HH11	1:A:608:ARG:CG	2.04	0.45
1:A:801:GLU:HG2	4:A:901:XB6:O5	2.17	0.45
1:A:231:PHE:HE1	1:A:249:VAL:HG12	1.81	0.45
1:A:231:PHE:CD1	1:A:231:PHE:O	2.70	0.45
2:B:368:GLU:H	2:B:369:PRO:HD3	1.81	0.45
2:B:397:LYS:HD3	2:B:398:TYR:CE1	2.52	0.45
2:B:437:TRP:O	2:B:437:TRP:CE3	2.70	0.45
1:A:721:ASP:O	1:A:725:GLY:N	2.46	0.45
2:B:412:LYS:HE3	2:B:412:LYS:HA	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:568:ARG:HH11	1:A:698:TYR:C	2.22	0.44
2:B:361:GLU:C	2:B:363:LEU:N	2.72	0.44
1:A:527:GLN:O	1:A:530:ASP:HB2	2.17	0.44
1:A:530:ASP:OD2	1:A:685:THR:HA	2.17	0.44
1:A:568:ARG:HD2	1:A:697:LEU:O	2.17	0.44
1:A:660:ASN:ND2	1:A:716:GLU:OE1	2.47	0.44
2:B:440:GLU:CD	2:B:440:GLU:C	2.85	0.44
1:A:317:VAL:HG13	1:A:571:TYR:HB3	1.99	0.44
1:A:191:GLN:HG2	1:A:255:TYR:OH	2.17	0.44
2:B:424:TYR:HA	2:B:427:ARG:HG3	1.99	0.44
1:A:793:ILE:CG2	1:A:794:PRO:CD	2.92	0.44
1:A:494:TYR:CD1	1:A:494:TYR:O	2.70	0.43
2:B:405:ILE:O	2:B:409:ILE:CD1	2.66	0.43
1:A:506:GLU:OE1	1:A:506:GLU:CA	2.66	0.43
1:A:496:GLU:O	1:A:499:GLU:HB2	2.18	0.43
1:A:548:SER:O	1:A:552:TRP:HB3	2.19	0.43
2:B:409:ILE:H	2:B:409:ILE:CD1	2.30	0.43
1:A:280:LYS:HD2	1:A:303:ASP:O	2.19	0.43
2:B:395:ILE:HG22	2:B:433:VAL:HG12	1.98	0.43
1:A:353:LEU:N	1:A:353:LEU:CD2	2.81	0.43
1:A:214:ARG:HD2	1:A:218:LEU:HD11	2.01	0.43
1:A:458:LEU:HD23	1:A:458:LEU:HA	1.67	0.43
1:A:538:PHE:CE1	1:A:706:LEU:HD13	2.53	0.43
1:A:691:LEU:HD22	1:A:705:ALA:HB1	2.01	0.43
1:A:392:LEU:CD2	1:A:398:PHE:CD2	3.02	0.43
1:A:364:GLU:HG2	1:A:370:VAL:HG23	1.99	0.43
2:B:355:THR:HG22	2:B:356:ASN:HD22	1.83	0.42
1:A:346:SER:CB	1:A:351:MET:HE3	2.49	0.42
1:A:677:LEU:HA	1:A:694:PHE:O	2.19	0.42
1:A:371:PRO:HG2	1:A:374:LYS:HB2	2.02	0.42
1:A:632:GLN:HG2	1:A:758:ARG:HD2	2.02	0.42
2:B:337:GLN:O	2:B:341:GLU:N	2.36	0.42
2:B:363:LEU:HD22	2:B:363:LEU:HA	1.84	0.42
1:A:670:PHE:CE2	1:A:740:VAL:HG22	2.54	0.42
1:A:657:GLY:O	1:A:752:ARG:NH2	2.52	0.42
1:A:724:VAL:HG11	1:A:746:THR:HG21	2.02	0.42
1:A:671:TRP:O	1:A:673:PRO:HD3	2.19	0.42
1:A:700:ALA:HB1	1:A:701:PRO:CD	2.41	0.42
1:A:235:LEU:HD13	1:A:235:LEU:HA	1.86	0.42
1:A:275:THR:O	1:A:275:THR:OG1	2.31	0.42
1:A:793:ILE:CG2	1:A:794:PRO:HD3	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:403:GLN:C	2:B:403:GLN:HE21	2.28	0.42
2:B:409:ILE:N	2:B:409:ILE:CD1	2.79	0.42
1:A:198:ASP:OD1	1:A:198:ASP:N	2.53	0.42
1:A:474:ILE:HD12	1:A:474:ILE:HA	1.95	0.42
1:A:322:LYS:O	1:A:325:TYR:HB2	2.19	0.42
1:A:640:VAL:HA	1:A:641:PRO:HA	1.90	0.42
4:A:901:XB6:C24	4:A:901:XB6:C11	2.98	0.42
1:A:208:LYS:HB3	1:A:208:LYS:HE2	1.74	0.41
1:A:715:MET:HE1	1:A:726:ARG:NH1	2.34	0.41
1:A:495:ASP:OD2	2:B:371:ARG:NH1	2.53	0.41
2:B:320:ASP:O	2:B:324:VAL:HG23	2.21	0.41
1:A:524:ARG:NH2	1:A:524:ARG:HG3	2.35	0.41
1:A:418:LEU:HD23	1:A:418:LEU:HA	1.72	0.41
1:A:691:LEU:HD13	1:A:727:CYS:SG	2.60	0.41
1:A:693:LEU:HD12	1:A:694:PHE:H	1.86	0.41
1:A:752:ARG:O	1:A:752:ARG:HG3	2.20	0.41
2:B:437:TRP:CE3	2:B:437:TRP:C	2.99	0.41
1:A:385:LEU:HD23	1:A:385:LEU:HA	1.81	0.41
1:A:683:SER:OG	1:A:684:THR:HG23	2.20	0.41
1:A:786:ILE:HD12	1:A:786:ILE:H	1.86	0.41
1:A:731:LEU:HD13	1:A:731:LEU:HA	1.84	0.41
1:A:568:ARG:NH1	1:A:699:LYS:CA	2.83	0.41
1:A:592:GLN:OE1	1:A:603:ILE:HD12	2.20	0.41
1:A:793:ILE:HG23	1:A:828:GLN:OE1	2.20	0.41
1:A:454:LYS:HB2	1:A:454:LYS:HE2	1.90	0.41
1:A:470:PRO:HA	1:A:471:PRO:C	2.45	0.41
1:A:214:ARG:HD2	1:A:218:LEU:CD1	2.51	0.40
4:A:901:XB6:O3	4:A:901:XB6:C9	2.69	0.40
1:A:520:TYR:CD1	1:A:520:TYR:C	2.96	0.40
2:B:429:ASN:N	2:B:429:ASN:ND2	2.69	0.40
1:A:221:TRP:CD1	1:A:262:ILE:HA	2.57	0.40
1:A:352:GLU:HG2	1:A:568:ARG:HB3	2.03	0.40
1:A:670:PHE:CZ	1:A:740:VAL:HG22	2.56	0.40
1:A:508:LEU:HD23	1:A:508:LEU:H	1.86	0.40
2:B:337:GLN:OE1	2:B:337:GLN:CA	2.69	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%)	630 (95%)	34 (5%)	0	100	100
2	B	131/144 (91%)	119 (91%)	10 (8%)	2 (2%)	8	28
3	C	2/9 (22%)	0	2 (100%)	0	100	100
All	All	797/1024 (78%)	749 (94%)	46 (6%)	2 (0%)	36	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	399	GLY
2	B	402	PHE

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%)	528 (93%)	38 (7%)	15	42
2	B	117/125 (94%)	92 (79%)	25 (21%)	1	4
3	C	4/9 (44%)	4 (100%)	0	100	100
All	All	687/849 (81%)	624 (91%)	63 (9%)	8	27

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	187	ARG

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Mol	Chain	Res	Type
1	A	199	ILE
1	A	200	ILE
1	A	235	LEU
1	A	273	LEU
1	A	286	SER
1	A	351	MET
1	A	370	VAL
1	A	374	LYS
1	A	377	MET
1	A	456	LYS
1	A	458	LEU
1	A	486	ASP
1	A	490	LEU
1	A	492	LYS
1	A	499	GLU
1	A	503	LYS
1	A	506	GLU
1	A	508	LEU
1	A	511	LEU
1	A	526	ARG
1	A	568	ARG
1	A	569	ASN
1	A	591	ARG
1	A	592	GLN
1	A	608	ARG
1	A	626	PRO
1	A	640	VAL
1	A	645	GLU
1	A	696	ASN
1	A	704	LEU
1	A	724	VAL
1	A	730	ILE
1	A	731	LEU
1	A	744	LYS
1	A	748	VAL
1	A	750	ARG
1	A	793	ILE
2	B	334	VAL
2	B	338	LEU
2	B	347	ARG
2	B	350	GLN
2	B	354	GLN

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Mol	Chain	Res	Type
2	B	355	THR
2	B	363	LEU
2	B	364	ASP
2	B	368	GLU
2	B	370	TYR
2	B	376	ILE
2	B	377	GLN
2	B	379	CYS
2	B	386	GLU
2	B	402	PHE
2	B	403	GLN
2	B	405	ILE
2	B	409	ILE
2	B	412	LYS
2	B	415	VAL
2	B	416	GLN
2	B	420	PHE
2	B	427	ARG
2	B	429	ASN
2	B	430	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	185	HIS
1	A	191	GLN
1	A	205	GLN
1	A	207	GLN
1	A	228	GLN
1	A	236	GLN
1	A	296	GLN
1	A	380	GLN
1	A	403	ASN
1	A	422	HIS
1	A	446	ASN
1	A	460	GLN
1	A	484	HIS
1	A	638	GLN
1	A	680	HIS
2	B	350	GLN
2	B	354	GLN
2	B	393	GLN

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Mol	Chain	Res	Type
2	B	403	GLN
2	B	423	ASN
2	B	429	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	$\# Z > 2$
4	XB6	A	901	-	75,76,76	1.38	8 (10%)	100,114,114	0.85	3 (3%)

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
4	XB6	A	901	-	-	20/47/63/63	0/8/8/8

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	XB6	C5-N1	-6.47	1.34	1.42
4	A	901	XB6	C24-C27	-3.71	1.34	1.43
4	A	901	XB6	P2-O10	-3.34	1.55	1.59
4	A	901	XB6	O4-C29	-3.14	1.36	1.43
4	A	901	XB6	P1-O10	-2.75	1.56	1.59
4	A	901	XB6	C6-C7	-2.54	1.36	1.39
4	A	901	XB6	C28-C29	-2.19	1.49	1.52
4	A	901	XB6	O5-C30	-2.03	1.37	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	XB6	C24-C27-N3	3.56	117.20	110.94
4	A	901	XB6	O11-P2-O12	2.28	123.06	112.44
4	A	901	XB6	C6-C5-N1	-2.15	119.48	122.70

There are no chirality outliers.

All (20) torsion outliers are listed below:

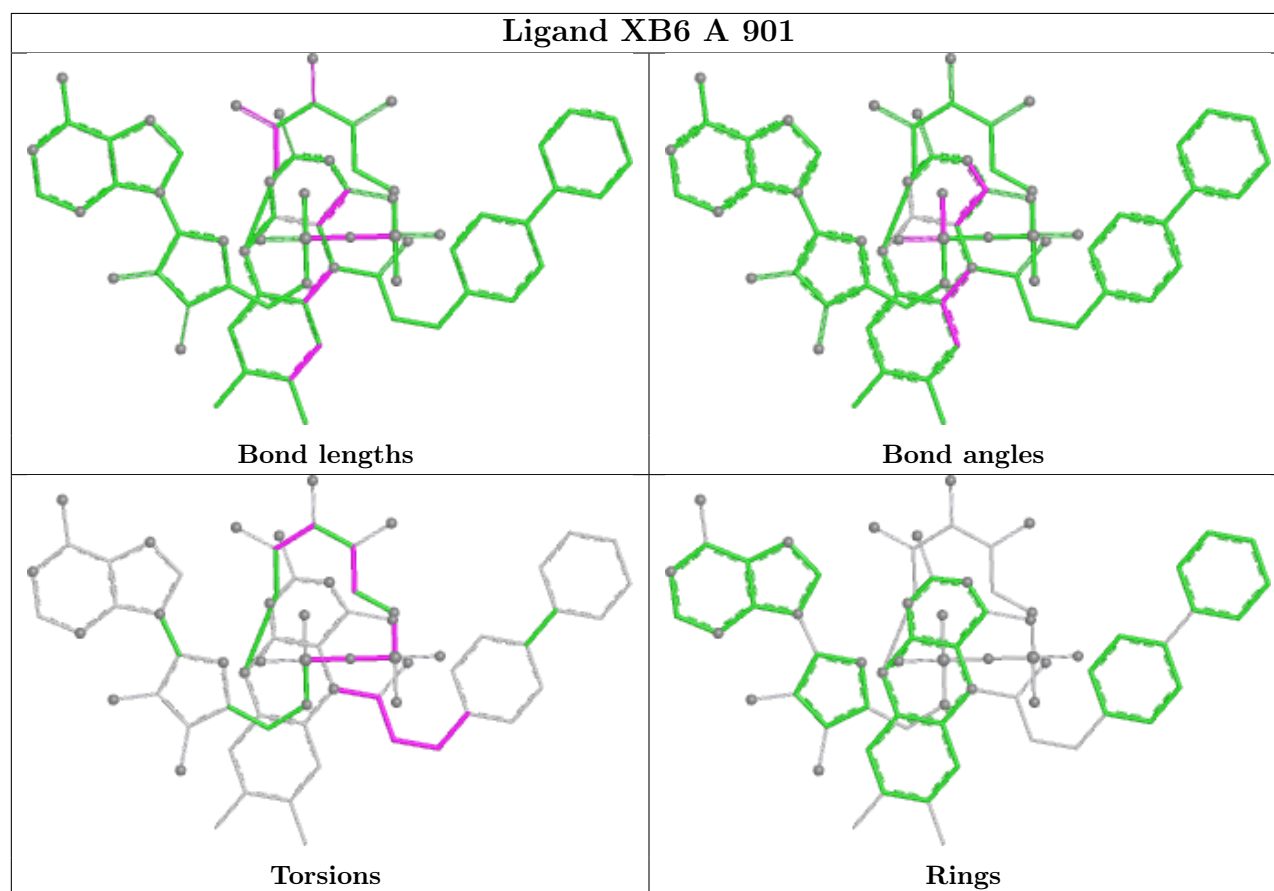
Mol	Chain	Res	Type	Atoms
4	A	901	XB6	C9-C10-C11-C12
4	A	901	XB6	C10-C9-N1-C5
4	A	901	XB6	C10-C9-N1-C24
4	A	901	XB6	O1-C9-N1-C5
4	A	901	XB6	O1-C9-N1-C24
4	A	901	XB6	C30-C31-C32-O7
4	A	901	XB6	O6-C31-C32-O7
4	A	901	XB6	C32-O7-P1-O10
4	A	901	XB6	C32-O7-P1-O8
4	A	901	XB6	C11-C10-C9-N1
4	A	901	XB6	O4-C29-C30-C31
4	A	901	XB6	O4-C29-C30-O5
4	A	901	XB6	C11-C10-C9-O1
4	A	901	XB6	P2-O10-P1-O7
4	A	901	XB6	P1-O10-P2-O12
4	A	901	XB6	C32-O7-P1-O9
4	A	901	XB6	P1-O10-P2-O11
4	A	901	XB6	C10-C11-C12-C23
4	A	901	XB6	C10-C11-C12-C13
4	A	901	XB6	C28-C29-C30-O5

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	901	XB6	3	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 ⓘ

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

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	666/871 (76%)	0.24	20 (3%) 52 42	62, 95, 128, 144	0
2	B	133/144 (92%)	0.73	9 (6%) 23 17	91, 123, 141, 154	0
3	C	4/9 (44%)	2.97	4 (100%) 0 0	108, 111, 118, 119	0
All	All	803/1024 (78%)	0.33	33 (4%) 41 33	62, 102, 133, 154	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	836	LEU	7.8
1	A	555	ASP	5.0
2	B	383	TRP	3.7
3	C	66	PRO	3.6
1	A	171	PRO	3.4
1	A	358	GLN	3.2
2	B	399	GLY	3.2
1	A	556	ASP	3.1
3	C	68	SER	3.1
1	A	833	MET	2.9
1	A	554	GLN	2.9
2	B	376	ILE	2.8
2	B	439	ALA	2.8
1	A	507	LYS	2.7
3	C	67	ARG	2.7
2	B	382	ARG	2.7
1	A	373	GLU	2.6
1	A	786	ILE	2.6
1	A	453	GLU	2.5
1	A	834	TYR	2.5
3	C	69	PHE	2.5
1	A	511	LEU	2.4
2	B	440	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	611	SER	2.3
2	B	308	ARG	2.3
1	A	518	ASP	2.2
1	A	503	LYS	2.2
1	A	272	PRO	2.2
1	A	363	TYR	2.2
2	B	386	GLU	2.2
1	A	377	MET	2.1
1	A	667	ASP	2.1
2	B	418	LYS	2.1

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](#)

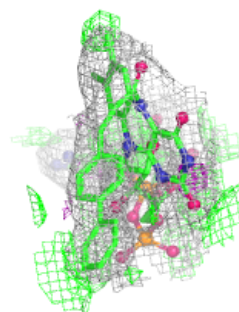
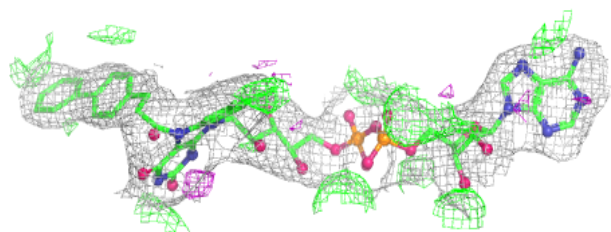
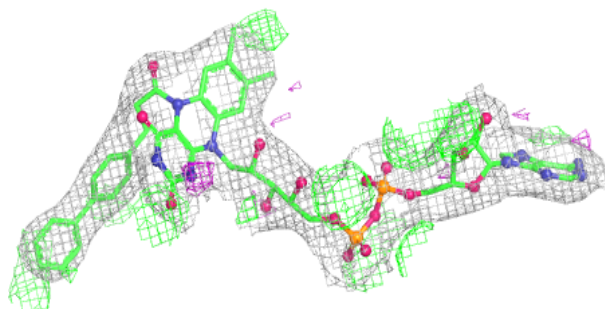
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	XB6	A	901	69/69	0.97	0.09	54,74,102,107	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.

Electron density around XB6 A 901:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.