



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:16 PM UTC

PDB ID : 8FJ4 / pdb_00008fj4
Title : LSD1-CoREST in complex with T108, short soaking
Authors : Caroli, J.; Mattevi, A.
Deposited on : 2022-12-19
Resolution : 2.76 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

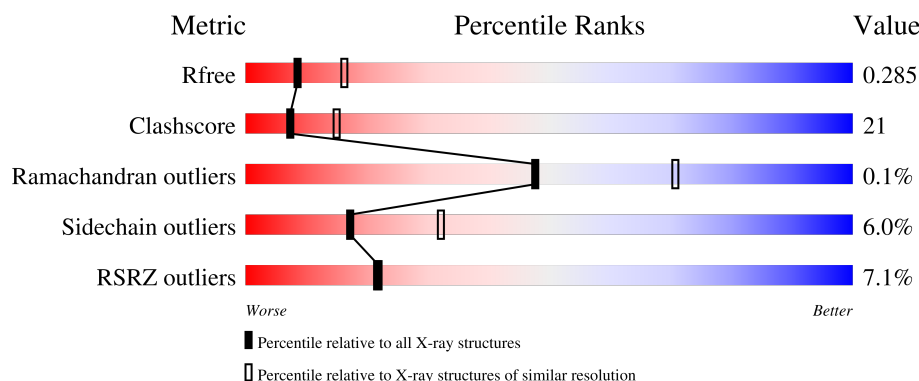
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.76 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	4% 53% 22% • 24%
2	B	144	14% 51% 30% 11% 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6434 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

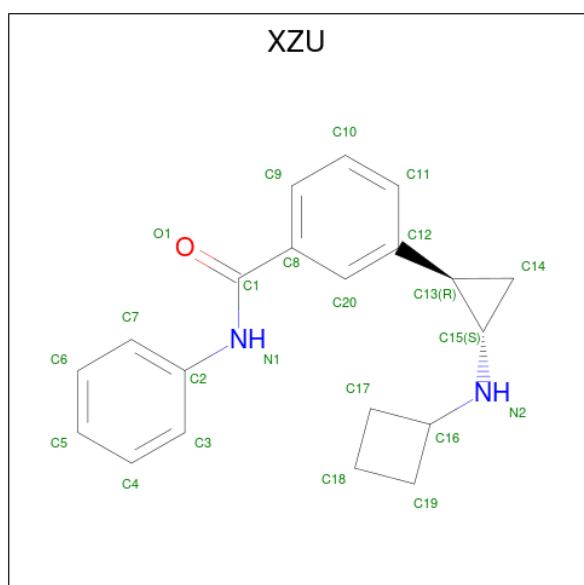
- Molecule 2 is a protein called REST corepressor 1.

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:

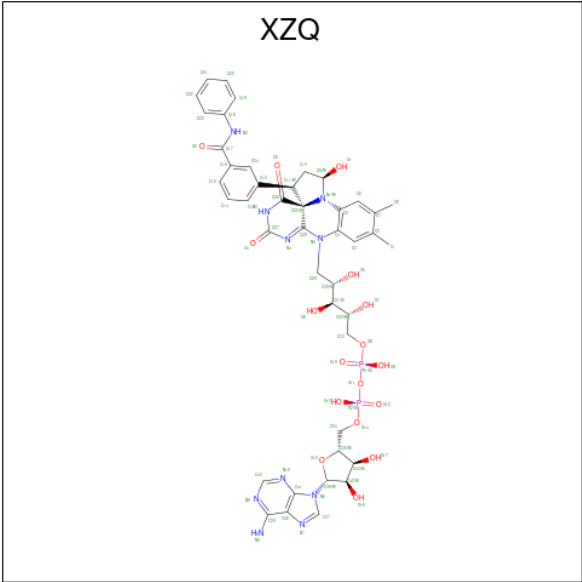
Chain	Residue	Modelled	Actual	Comment	Reference
B	297	GLY	-	expression tag	UNP Q9UKL0
B	298	PRO	-	expression tag	UNP Q9UKL0
B	299	LEU	-	expression tag	UNP Q9UKL0
B	300	GLY	-	expression tag	UNP Q9UKL0
B	301	SER	-	expression tag	UNP Q9UKL0
B	302	PRO	-	expression tag	UNP Q9UKL0
B	303	GLU	-	expression tag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

- Molecule 3 is 3-[(1R,2S)-2-(cyclobutylamino)cyclopropyl]-N-phenylbenzamide (CCD ID: XZU) (formula: C₂₀H₂₂N₂O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			23	20	2	1		
3	A	1	Total	C	N	O	0	0
			23	20	2	1		
3	A	1	Total	C	N	O	0	0
			23	20	2	1		

- Molecule 4 is [(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-3,4-dihydroxyoxolan-2-yl]methyl (2R,3S,4S)-2,3,4-trihydroxy-5-[(1R,3R,3aS,13R)-1-hydroxy-10,11-dimethyl-4,6-dioxo-3-[3-(phenylcarbamoyl)phenyl]-2,3,5,6-tetrahydro-1H-benzo[g]pyrrolo[2,1-e]pteridin-8(4H)-yl]pentyl dihydrogen diphosphate (CCD ID: XZQ) (formula: C₄₃H₄₈N₁₀O₁₇P₂) (labeled as "Ligand of Interest" by depositor).

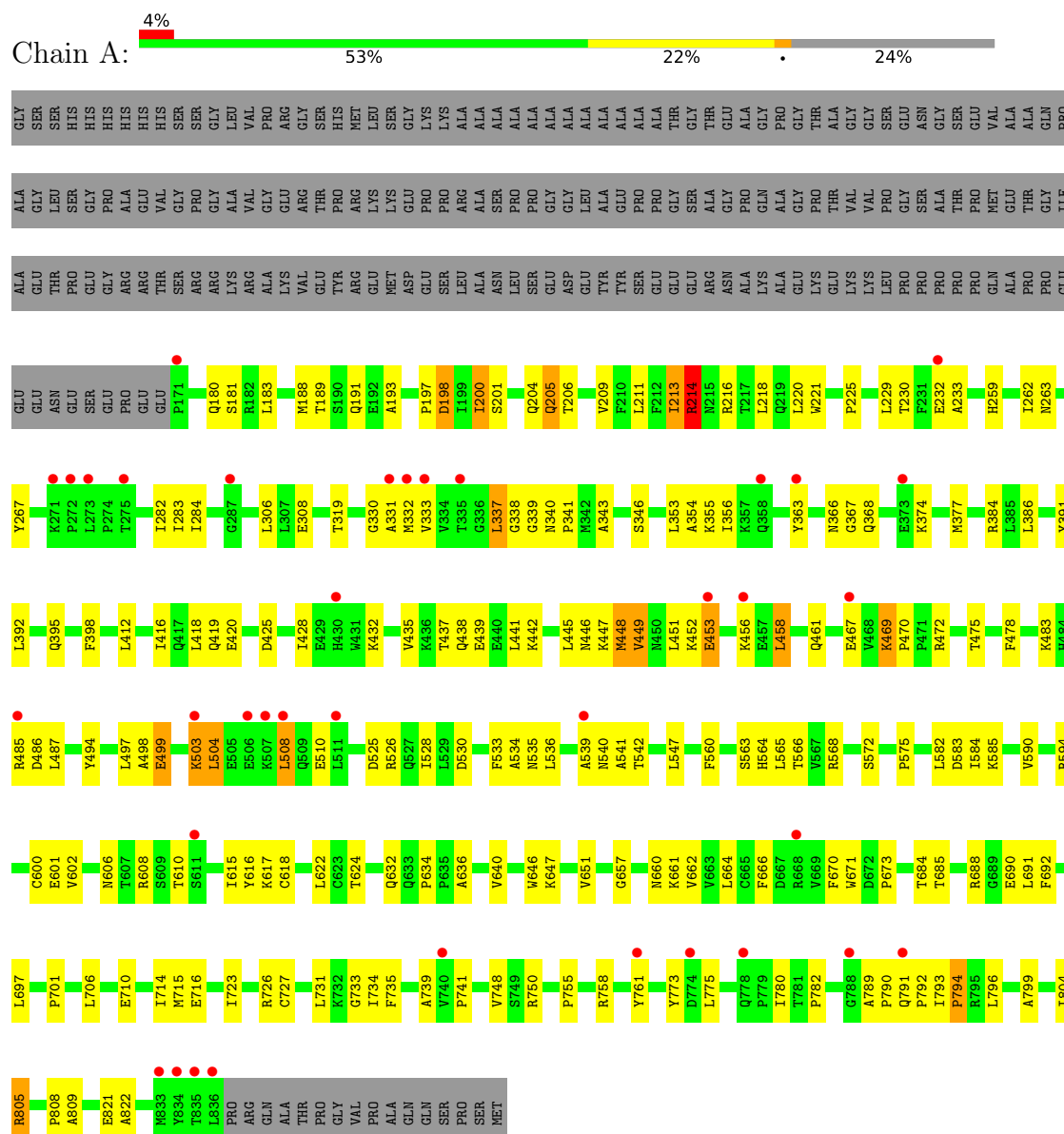


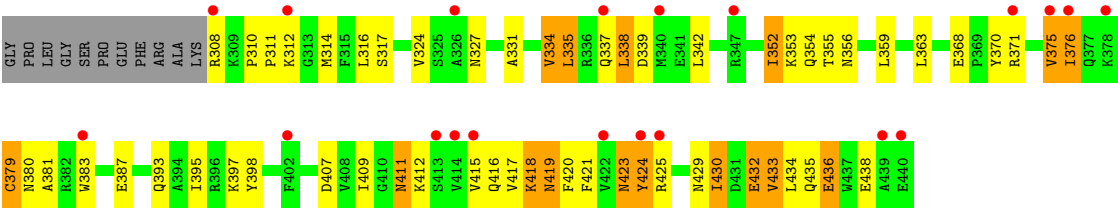
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	A	1	72	43	10	17	2	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 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	118.35Å 179.74Å 232.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.78 – 2.76 48.78 – 2.76	Depositor EDS
% Data completeness (in resolution range)	97.8 (48.78-2.76) 97.8 (48.78-2.76)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.18.2	Depositor
R, R_{free}	0.224 , 0.248 0.265 , 0.285	Depositor DCC
R_{free} test set	2000 reflections (3.14%)	wwPDB-VP
Wilson B-factor (Å ²)	80.6	Xtriage
Anisotropy	0.461	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 85.7	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	0.94	EDS
Total number of atoms	6434	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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: XZU, XZQ

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	0.71	4/5331 (0.1%)	0.92	4/7232 (0.1%)
2	B	0.77	0/1091	1.23	2/1471 (0.1%)
All	All	0.72	4/6422 (0.1%)	0.98	6/8703 (0.1%)

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	470	PRO	C-O	-6.44	1.19	1.24
1	A	341	PRO	C-O	-5.54	1.17	1.24
1	A	794	PRO	C-O	-5.47	1.18	1.23
1	A	808	PRO	C-O	-5.07	1.17	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	499	GLU	N-CA-C	-9.17	100.85	111.03
2	B	417	VAL	N-CA-C	-9.13	101.50	110.72
2	B	411	ASN	N-CA-C	-7.20	105.03	113.88
1	A	214	ARG	CA-C-O	-5.20	115.36	120.82
1	A	469	LYS	CA-C-N	5.06	123.31	119.66
1	A	469	LYS	C-N-CA	5.06	123.31	119.66

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	214	ARG	Mainchain

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	212	0
2	B	1076	0	1091	90	0
3	A	69	0	0	2	0
4	A	72	0	0	5	0
All	All	6434	0	6343	262	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 21.

All (262) 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)
2:B:419:ASN:C	2:B:423:ASN:HD21	1.33	1.37
2:B:419:ASN:O	2:B:423:ASN:ND2	1.72	1.20
1:A:441:LEU:O	1:A:445:LEU:HD12	1.48	1.13
1:A:188:MET:HE3	1:A:200:ILE:HG13	1.27	1.11
1:A:448:MET:CE	1:A:497:LEU:HB3	1.85	1.05
1:A:449:VAL:HG23	2:B:363:LEU:CD1	1.86	1.04
1:A:188:MET:CE	1:A:200:ILE:HG13	1.87	1.04
1:A:441:LEU:O	1:A:445:LEU:CD1	2.07	1.01
1:A:448:MET:HE2	1:A:497:LEU:HB3	1.38	1.01
2:B:419:ASN:C	2:B:423:ASN:ND2	2.16	0.98
1:A:449:VAL:HG23	2:B:363:LEU:HD11	1.46	0.97
1:A:209:VAL:O	1:A:213:ILE:HG13	1.65	0.97
1:A:374:LYS:NZ	1:A:525:ASP:OD1	1.98	0.96
1:A:392:LEU:HD11	2:B:316:LEU:CD2	1.99	0.93
1:A:445:LEU:HD12	1:A:445:LEU:H	1.30	0.92
1:A:449:VAL:CG2	2:B:363:LEU:CD1	2.49	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:416:GLN:HA	2:B:419:ASN:HB2	1.55	0.89
2:B:420:PHE:C	2:B:423:ASN:HD22	1.81	0.88
2:B:420:PHE:CA	2:B:423:ASN:ND2	2.37	0.87
2:B:387:GLU:OE1	2:B:411:ASN:ND2	2.07	0.87
2:B:425:ARG:HA	2:B:430:ILE:HG13	1.58	0.86
1:A:188:MET:HE1	1:A:200:ILE:CB	2.06	0.85
1:A:449:VAL:CG2	2:B:363:LEU:HD11	2.05	0.85
2:B:387:GLU:CD	2:B:411:ASN:HD21	1.84	0.84
1:A:188:MET:CE	1:A:200:ILE:CG1	2.56	0.83
1:A:188:MET:HE1	1:A:200:ILE:HB	1.61	0.81
1:A:566:THR:HG21	1:A:697:LEU:HD22	1.63	0.80
2:B:324:VAL:HG13	2:B:331:ALA:HB2	1.63	0.80
1:A:437:THR:OG1	1:A:508:LEU:HD21	1.81	0.80
1:A:441:LEU:HG	1:A:445:LEU:HD11	1.65	0.79
1:A:448:MET:HE3	1:A:497:LEU:HB3	1.67	0.77
2:B:423:ASN:ND2	2:B:423:ASN:H	1.83	0.77
1:A:188:MET:HE3	1:A:200:ILE:CG1	2.09	0.77
1:A:486:ASP:OD2	2:B:398:TYR:OH	2.01	0.77
1:A:197:PRO:O	1:A:201:SER:HB3	1.86	0.75
1:A:188:MET:CE	1:A:200:ILE:CB	2.64	0.75
1:A:461:GLN:OE1	1:A:483:LYS:NZ	2.13	0.75
4:A:1404:XZQ:O3	4:A:1404:XZQ:O1	2.05	0.74
1:A:448:MET:HE2	1:A:497:LEU:CB	2.16	0.74
1:A:441:LEU:C	1:A:445:LEU:CD1	2.61	0.74
1:A:793:ILE:HD12	1:A:793:ILE:H	1.52	0.73
1:A:437:THR:OG1	1:A:508:LEU:CD2	2.37	0.72
1:A:188:MET:HE1	1:A:200:ILE:CA	2.19	0.72
1:A:446:ASN:OD1	2:B:359:LEU:HD21	1.89	0.72
2:B:311:PRO:HG2	2:B:314:MET:HG3	1.71	0.72
1:A:230:THR:OG1	1:A:232:GLU:OE1	2.06	0.72
1:A:392:LEU:HD11	2:B:316:LEU:HD22	1.70	0.72
1:A:449:VAL:HG23	2:B:363:LEU:HD13	1.69	0.72
1:A:439:GLU:HG2	2:B:352:ILE:CD1	2.20	0.72
1:A:384:ARG:NH2	2:B:312:LYS:O	2.23	0.71
2:B:420:PHE:C	2:B:423:ASN:ND2	2.48	0.71
1:A:821:GLU:HA	1:A:821:GLU:OE1	1.91	0.71
1:A:453:GLU:OE1	1:A:453:GLU:HA	1.89	0.71
1:A:503:LYS:HG2	1:A:504:LEU:HD23	1.72	0.71
1:A:392:LEU:CD1	2:B:316:LEU:HD22	2.21	0.70
1:A:439:GLU:CG	2:B:352:ILE:CD1	2.70	0.70
1:A:690:GLU:OE2	1:A:726:ARG:NH1	2.24	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:445:LEU:HD12	1:A:445:LEU:N	2.06	0.69
1:A:230:THR:HG23	1:A:233:ALA:H	1.56	0.69
2:B:420:PHE:N	2:B:423:ASN:HD21	1.88	0.69
1:A:448:MET:CE	1:A:497:LEU:C	2.65	0.69
1:A:564:HIS:ND1	4:A:1404:XZQ:C19	2.55	0.69
1:A:425:ASP:OD2	2:B:338:LEU:CD1	2.41	0.69
1:A:188:MET:CE	1:A:200:ILE:HB	2.22	0.68
2:B:420:PHE:HA	2:B:423:ASN:ND2	2.07	0.68
1:A:458:LEU:HB3	1:A:487:LEU:HD12	1.78	0.66
1:A:438:GLN:CD	1:A:508:LEU:HD11	2.21	0.66
2:B:395:ILE:HG22	2:B:433:VAL:CG1	2.26	0.66
2:B:376:ILE:HG12	2:B:376:ILE:O	1.96	0.65
1:A:337:LEU:HD23	1:A:337:LEU:N	2.10	0.65
2:B:420:PHE:N	2:B:423:ASN:ND2	2.45	0.64
2:B:436:GLU:OE1	2:B:436:GLU:HA	1.96	0.64
1:A:526:ARG:NH1	1:A:530:ASP:OD1	2.31	0.63
1:A:606:ASN:HD21	1:A:608:ARG:HE	1.44	0.63
1:A:395:GLN:HE21	2:B:308:ARG:N	1.96	0.63
1:A:499:GLU:HA	1:A:499:GLU:OE1	1.98	0.62
1:A:209:VAL:O	1:A:213:ILE:CG1	2.42	0.62
1:A:330:GLY:O	1:A:661:LYS:NZ	2.33	0.62
1:A:439:GLU:HG2	2:B:352:ILE:HD13	1.81	0.62
1:A:209:VAL:C	1:A:213:ILE:HG13	2.25	0.62
1:A:206:THR:O	1:A:209:VAL:HB	2.00	0.62
2:B:335:LEU:N	2:B:335:LEU:HD23	2.15	0.61
1:A:671:TRP:HA	1:A:735:PHE:CE2	2.35	0.61
1:A:198:ASP:OD1	1:A:198:ASP:N	2.22	0.61
1:A:331:ALA:HA	4:A:1404:XZQ:N1	2.16	0.61
1:A:353:LEU:HB3	1:A:565:LEU:HD23	1.83	0.60
1:A:425:ASP:OD2	2:B:338:LEU:HD13	2.01	0.60
2:B:375:VAL:HG12	2:B:375:VAL:O	2.00	0.60
2:B:338:LEU:N	2:B:338:LEU:HD23	2.16	0.60
1:A:789:ALA:HB1	1:A:790:PRO:HD2	1.84	0.60
1:A:804:ILE:HG23	1:A:804:ILE:O	2.02	0.59
1:A:432:LYS:HA	1:A:435:VAL:HG22	1.85	0.59
1:A:467:GLU:OE2	1:A:467:GLU:HA	2.04	0.58
1:A:283:ILE:HG12	1:A:622:LEU:HB3	1.85	0.58
1:A:445:LEU:CD1	1:A:445:LEU:H	2.08	0.58
1:A:780:ILE:HB	1:A:796:LEU:HB3	1.85	0.58
1:A:755:PRO:HA	1:A:758:ARG:NE	2.19	0.58
1:A:438:GLN:CD	1:A:508:LEU:CD1	2.76	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:439:GLU:HG3	2:B:352:ILE:CD1	2.34	0.57
1:A:438:GLN:HG2	1:A:508:LEU:HD11	1.85	0.57
1:A:449:VAL:HG22	2:B:363:LEU:CD1	2.34	0.57
1:A:441:LEU:HD23	2:B:356:ASN:HD22	1.69	0.57
1:A:308:GLU:OE1	4:A:1404:XZQ:O17	2.23	0.57
1:A:333:VAL:HG21	4:A:1404:XZQ:C17	2.35	0.56
1:A:448:MET:CE	1:A:497:LEU:CB	2.74	0.56
2:B:434:LEU:O	2:B:438:GLU:HG3	2.05	0.56
1:A:535:ASN:OD1	1:A:692:PHE:HE1	1.88	0.56
1:A:755:PRO:HA	1:A:758:ARG:HE	1.69	0.56
1:A:671:TRP:HA	1:A:735:PHE:HE2	1.70	0.56
1:A:448:MET:HE1	1:A:497:LEU:C	2.31	0.56
1:A:448:MET:HE1	1:A:498:ALA:HA	1.88	0.55
1:A:478:PHE:CD1	2:B:393:GLN:HB3	2.41	0.55
1:A:666:PHE:O	1:A:701:PRO:HG2	2.05	0.55
1:A:448:MET:HB3	2:B:363:LEU:HD21	1.89	0.55
1:A:475:THR:HA	2:B:393:GLN:HE22	1.72	0.55
1:A:442:LYS:HD2	2:B:355:THR:HG21	1.89	0.54
1:A:448:MET:HE1	1:A:498:ALA:CA	2.37	0.54
1:A:284:ILE:HG12	1:A:590:VAL:HG21	1.88	0.54
1:A:441:LEU:CG	1:A:445:LEU:HD11	2.35	0.54
1:A:671:TRP:O	1:A:673:PRO:HD3	2.08	0.53
1:A:340:ASN:HB2	1:A:560:PHE:CD1	2.44	0.53
1:A:541:ALA:O	1:A:657:GLY:HA3	2.09	0.53
1:A:438:GLN:HE22	2:B:353:LYS:HG3	1.73	0.53
2:B:416:GLN:OE1	2:B:416:GLN:N	2.30	0.53
1:A:438:GLN:CG	1:A:508:LEU:HD11	2.39	0.53
1:A:205:GLN:O	1:A:209:VAL:HG23	2.09	0.53
1:A:343:ALA:O	1:A:346:SER:OG	2.25	0.53
1:A:391:TYR:CZ	2:B:310:PRO:HD3	2.44	0.52
1:A:229:LEU:N	1:A:263:ASN:OD1	2.37	0.52
1:A:209:VAL:HG12	1:A:213:ILE:HD11	1.90	0.52
2:B:424:TYR:N	2:B:424:TYR:CD2	2.78	0.52
2:B:383:TRP:CD2	2:B:412:LYS:HE2	2.45	0.52
1:A:526:ARG:HH11	1:A:526:ARG:HG3	1.75	0.51
1:A:395:GLN:NE2	2:B:308:ARG:N	2.57	0.51
2:B:423:ASN:HB2	2:B:424:TYR:CD2	2.45	0.51
1:A:439:GLU:HG3	2:B:352:ILE:HD12	1.92	0.51
2:B:379:CYS:SG	2:B:380:ASN:N	2.84	0.51
2:B:420:PHE:CA	2:B:423:ASN:HD22	2.11	0.51
2:B:415:VAL:O	2:B:418:LYS:HB2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:383:TRP:CH2	2:B:412:LYS:HG3	2.45	0.50
1:A:448:MET:HE1	1:A:498:ALA:N	2.26	0.50
2:B:423:ASN:HB2	2:B:424:TYR:HD2	1.77	0.50
1:A:363:TYR:CD2	1:A:734:ILE:HG23	2.46	0.50
1:A:366:ASN:OD1	1:A:367:GLY:N	2.44	0.50
1:A:773:TYR:HB2	1:A:805:ARG:HB2	1.93	0.50
2:B:334:VAL:HG12	2:B:335:LEU:HD23	1.94	0.49
1:A:354:ALA:HB2	1:A:568:ARG:HD2	1.93	0.49
1:A:216:ARG:HG3	1:A:220:LEU:CD1	2.43	0.49
1:A:716:GLU:OE2	1:A:750:ARG:HB3	2.11	0.49
1:A:282:ILE:HG21	1:A:602:VAL:HG21	1.94	0.49
1:A:504:LEU:HD23	1:A:504:LEU:N	2.28	0.49
3:A:1403:XZU:C9	3:A:1403:XZU:C3	2.91	0.49
2:B:409:ILE:HG22	2:B:411:ASN:H	1.77	0.48
1:A:485:ARG:HD3	1:A:485:ARG:C	2.38	0.48
1:A:367:GLY:HA3	1:A:733:GLY:O	2.14	0.48
1:A:181:SER:OG	1:A:218:LEU:HD23	2.14	0.48
1:A:216:ARG:O	1:A:220:LEU:HD12	2.14	0.48
1:A:670:PHE:HZ	1:A:731:LEU:HD13	1.78	0.48
1:A:525:ASP:O	1:A:528:ILE:N	2.47	0.47
1:A:582:LEU:HB2	1:A:584:ILE:HD11	1.96	0.47
1:A:216:ARG:HG3	1:A:220:LEU:HD11	1.96	0.47
1:A:319:THR:HB	1:A:572:SER:HB3	1.96	0.47
1:A:340:ASN:CG	1:A:560:PHE:CE1	2.93	0.47
1:A:384:ARG:HB2	2:B:314:MET:HE3	1.95	0.47
1:A:366:ASN:OD1	1:A:368:GLN:N	2.45	0.47
1:A:456:LYS:HA	2:B:370:TYR:CE2	2.49	0.47
1:A:539:ALA:HA	3:A:1401:XZU:C10	2.45	0.47
1:A:793:ILE:HG22	1:A:794:PRO:CD	2.45	0.47
1:A:214:ARG:HG3	1:A:218:LEU:CD1	2.45	0.47
1:A:793:ILE:HG22	1:A:794:PRO:HD2	1.97	0.47
1:A:188:MET:HE1	1:A:200:ILE:HA	1.97	0.46
1:A:353:LEU:HD13	1:A:565:LEU:HD22	1.96	0.46
1:A:392:LEU:HD12	2:B:316:LEU:HD22	1.96	0.46
1:A:632:GLN:OE1	1:A:636:ALA:HB2	2.16	0.46
1:A:646:TRP:CZ3	1:A:647:LYS:HE3	2.51	0.46
2:B:395:ILE:HG22	2:B:433:VAL:HG12	1.96	0.46
2:B:421:PHE:HE1	2:B:434:LEU:HD11	1.81	0.46
1:A:451:LEU:HD23	1:A:494:TYR:HB2	1.97	0.46
1:A:540:ASN:HB3	1:A:547:LEU:HD21	1.98	0.46
1:A:624:THR:HG22	1:A:799:ALA:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:ILE:O	1:A:420:GLU:HG3	2.16	0.46
1:A:662:VAL:HG13	1:A:748:VAL:HG22	1.98	0.46
1:A:216:ARG:HA	1:A:216:ARG:HD2	1.80	0.46
1:A:338:GLY:O	1:A:560:PHE:HB3	2.16	0.46
1:A:664:LEU:HD11	1:A:727:CYS:SG	2.55	0.45
1:A:565:LEU:HD12	1:A:565:LEU:N	2.32	0.45
1:A:791:GLN:HA	1:A:792:PRO:HD3	1.81	0.45
1:A:386:LEU:HD23	1:A:386:LEU:HA	1.67	0.45
1:A:660:ASN:ND2	1:A:716:GLU:OE1	2.49	0.45
2:B:395:ILE:HG22	2:B:433:VAL:HG11	1.96	0.45
1:A:428:ILE:HG12	2:B:342:LEU:CD1	2.46	0.45
1:A:485:ARG:HG3	2:B:407:ASP:HB2	1.99	0.45
1:A:419:GLN:NE2	2:B:314:MET:HA	2.32	0.44
1:A:594:ARG:CZ	1:A:640:VAL:HG11	2.47	0.44
1:A:710:GLU:O	1:A:714:ILE:HG12	2.18	0.44
1:A:758:ARG:HA	1:A:758:ARG:HD3	1.80	0.44
1:A:425:ASP:OD2	2:B:338:LEU:HD11	2.15	0.44
1:A:600:CYS:HB2	1:A:618:CYS:SG	2.57	0.44
1:A:615:ILE:HG21	1:A:617:LYS:HE3	1.99	0.44
2:B:354:GLN:HE21	2:B:354:GLN:HB2	1.57	0.44
1:A:445:LEU:HB2	2:B:359:LEU:HD23	1.99	0.44
1:A:306:LEU:HD13	1:A:584:ILE:HG13	1.98	0.44
1:A:353:LEU:HB3	1:A:565:LEU:CD2	2.46	0.44
1:A:392:LEU:HD23	1:A:398:PHE:CD1	2.53	0.44
1:A:193:ALA:HB2	1:A:200:ILE:HD12	2.00	0.44
1:A:534:ALA:HB2	1:A:688:ARG:HG3	1.99	0.44
1:A:441:LEU:C	1:A:445:LEU:HD13	2.41	0.43
1:A:263:ASN:O	1:A:267:TYR:CE1	2.71	0.43
1:A:263:ASN:C	1:A:267:TYR:CE1	2.96	0.43
1:A:632:GLN:HG2	1:A:634:PRO:O	2.17	0.43
2:B:429:ASN:HB3	2:B:432:GLU:HB2	1.99	0.43
1:A:355:LYS:HD3	1:A:563:SER:CB	2.49	0.43
1:A:392:LEU:CD1	2:B:316:LEU:CD2	2.78	0.43
2:B:395:ILE:CG2	2:B:433:VAL:CG1	2.94	0.43
1:A:332:MET:SD	1:A:661:LYS:HE2	2.59	0.43
1:A:412:LEU:HD13	1:A:533:PHE:CE1	2.54	0.43
1:A:572:SER:O	1:A:575:PRO:HD2	2.18	0.43
1:A:438:GLN:NE2	2:B:353:LYS:HG3	2.34	0.43
1:A:583:ASP:OD1	1:A:585:LYS:NZ	2.51	0.43
1:A:647:LYS:O	1:A:651:VAL:HG23	2.19	0.43
1:A:478:PHE:CE1	2:B:393:GLN:HB3	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:469:LYS:HE2	1:A:469:LYS:HA	2.01	0.42
1:A:214:ARG:HG3	1:A:218:LEU:HD11	2.01	0.42
1:A:684:THR:HG22	1:A:685:THR:N	2.35	0.42
1:A:775:LEU:HD23	1:A:775:LEU:HA	1.79	0.42
2:B:368:GLU:OE1	2:B:371:ARG:NE	2.49	0.42
1:A:428:ILE:HG12	2:B:342:LEU:HD12	2.02	0.42
1:A:221:TRP:CD1	1:A:262:ILE:HA	2.55	0.42
1:A:183:LEU:HD23	1:A:189:THR:HG21	2.01	0.42
1:A:374:LYS:HE3	1:A:374:LYS:HB3	1.84	0.42
1:A:392:LEU:HD11	2:B:316:LEU:HD23	1.95	0.42
1:A:601:GLU:HA	1:A:616:TYR:O	2.20	0.42
1:A:356:ILE:HD11	1:A:566:THR:CG2	2.49	0.42
1:A:691:LEU:N	1:A:691:LEU:HD12	2.35	0.42
1:A:191:GLN:HE22	1:A:259:HIS:CD2	2.38	0.41
2:B:424:TYR:N	2:B:424:TYR:HD2	2.17	0.41
1:A:782:PRO:O	1:A:792:PRO:HG2	2.20	0.41
1:A:441:LEU:O	1:A:445:LEU:HD13	2.12	0.41
1:A:448:MET:HE3	1:A:497:LEU:C	2.44	0.41
2:B:379:CYS:HA	2:B:411:ASN:O	2.20	0.41
2:B:423:ASN:ND2	2:B:423:ASN:N	2.60	0.41
1:A:211:LEU:HD23	1:A:211:LEU:HA	1.94	0.41
1:A:333:VAL:HA	1:A:565:LEU:O	2.21	0.41
2:B:395:ILE:CG2	2:B:433:VAL:HG11	2.50	0.41
1:A:367:GLY:HA2	1:A:734:ILE:HG12	2.02	0.41
2:B:415:VAL:O	2:B:418:LYS:N	2.53	0.41
1:A:282:ILE:HD13	1:A:282:ILE:HA	1.93	0.40
1:A:715:MET:HE3	1:A:723:ILE:HG12	2.02	0.40
1:A:715:MET:CE	1:A:723:ILE:HG12	2.52	0.40
1:A:180:GLN:HA	1:A:339:GLY:HA2	2.04	0.40
1:A:374:LYS:O	1:A:377:MET:N	2.54	0.40
1:A:761:TYR:CD2	1:A:809:ALA:HB1	2.56	0.40
1:A:340:ASN:ND2	1:A:560:PHE:CE1	2.90	0.40
1:A:412:LEU:HD23	1:A:412:LEU:HA	1.95	0.40
1:A:536:LEU:HD12	1:A:536:LEU:HA	1.80	0.40
1:A:691:LEU:HA	1:A:706:LEU:O	2.21	0.40
1:A:739:ALA:O	1:A:741:PRO:HD3	2.21	0.40
1:A:821:GLU:O	1:A:822:ALA:C	2.64	0.40
1:A:418:LEU:HD12	2:B:324:VAL:HG21	2.03	0.40
2:B:380:ASN:OD1	2:B:381:ALA:N	2.55	0.40
2:B:419:ASN:HD22	2:B:419:ASN:HA	1.59	0.40
1:A:540:ASN:O	1:A:542:THR:HG22	2.21	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	664/871 (76%)	630 (95%)	33 (5%)	1 (0%)	43	64
2	B	131/144 (91%)	124 (95%)	7 (5%)	0	100	100
All	All	795/1015 (78%)	754 (95%)	40 (5%)	1 (0%)	48	71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	225	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	566/715 (79%)	546 (96%)	20 (4%)	32	56
2	B	117/125 (94%)	96 (82%)	21 (18%)	2	3
All	All	683/840 (81%)	642 (94%)	41 (6%)	17	32

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	198	ASP
1	A	200	ILE

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Mol	Chain	Res	Type
1	A	204	GLN
1	A	205	GLN
1	A	213	ILE
1	A	214	ARG
1	A	337	LEU
1	A	447	LYS
1	A	448	MET
1	A	449	VAL
1	A	452	LYS
1	A	453	GLU
1	A	458	LEU
1	A	472	ARG
1	A	503	LYS
1	A	504	LEU
1	A	508	LEU
1	A	510	GLU
1	A	610	THR
1	A	805	ARG
2	B	317	SER
2	B	327	ASN
2	B	334	VAL
2	B	335	LEU
2	B	337	GLN
2	B	338	LEU
2	B	339	ASP
2	B	352	ILE
2	B	375	VAL
2	B	376	ILE
2	B	379	CYS
2	B	397	LYS
2	B	418	LYS
2	B	419	ASN
2	B	423	ASN
2	B	424	TYR
2	B	430	ILE
2	B	432	GLU
2	B	433	VAL
2	B	435	GLN
2	B	436	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	GLN
1	A	191	GLN
1	A	204	GLN
1	A	250	HIS
1	A	259	HIS
1	A	380	GLN
1	A	410	GLN
1	A	430	HIS
1	A	554	GLN
1	A	812	HIS
2	B	351	ASN
2	B	354	GLN
2	B	419	ASN
2	B	423	ASN
2	B	435	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	XZU	A	1401	-	25,26,26	0.71	0	27,36,36	2.49	5 (18%)
4	XZQ	A	1404	-	77,80,80	1.32	11 (14%)	105,123,123	1.03	6 (5%)
3	XZU	A	1403	-	25,26,26	0.59	0	27,36,36	0.86	1 (3%)
3	XZU	A	1402	-	25,26,26	0.77	1 (4%)	27,36,36	1.36	3 (11%)

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	XZU	A	1401	-	-	4/12/27/27	0/4/4/4
4	XZQ	A	1404	-	-	12/46/114/114	0/8/9/9
3	XZU	A	1403	-	-	4/12/27/27	0/4/4/4
3	XZU	A	1402	-	-	0/12/27/27	0/4/4/4

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1404	XZQ	P2-O11	-3.50	1.55	1.59
4	A	1404	XZQ	C5-N1	-3.22	1.35	1.41
4	A	1404	XZQ	C27-N4	-3.15	1.29	1.36
4	A	1404	XZQ	P1-O11	-3.04	1.56	1.59
4	A	1404	XZQ	O5-C30	-2.85	1.37	1.43
4	A	1404	XZQ	C26-N3	-2.82	1.32	1.37
4	A	1404	XZQ	C5-C4	-2.18	1.37	1.41
4	A	1404	XZQ	C6-C7	-2.16	1.36	1.39
3	A	1402	XZU	C15-N2	-2.16	1.43	1.47
4	A	1404	XZQ	C9-N1	-2.11	1.43	1.46
4	A	1404	XZQ	O4-C27	-2.10	1.20	1.24
4	A	1404	XZQ	O6-C31	-2.05	1.37	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1401	XZU	C18-C17-C16	-6.94	84.67	88.29
3	A	1401	XZU	C18-C19-C16	-6.74	84.77	88.29
3	A	1401	XZU	C19-C16-C17	-6.20	84.59	88.51
4	A	1404	XZQ	C5-N1-C9	-5.10	112.21	123.92
3	A	1402	XZU	C18-C19-C16	-4.47	85.96	88.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1402	XZU	C18-C17-C16	-4.44	85.97	88.29
4	A	1404	XZQ	C26-C25-C28	-3.66	106.91	113.40
3	A	1401	XZU	C11-C12-C13	-3.40	114.56	121.08
3	A	1401	XZU	C20-C12-C13	3.10	126.11	120.11
4	A	1404	XZQ	C26-N3-C27	-2.81	121.19	125.42
4	A	1404	XZQ	C25-N1-C9	-2.40	107.09	109.50
4	A	1404	XZQ	O11-P1-O10	-2.24	103.97	110.70
3	A	1402	XZU	C19-C16-C17	-2.23	87.10	88.51
4	A	1404	XZQ	O13-P2-O12	2.22	122.78	112.44
3	A	1403	XZU	C2-N1-C1	2.02	131.91	126.61

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1401	XZU	C8-C1-N1-C2
3	A	1403	XZU	C8-C1-N1-C2
3	A	1403	XZU	O1-C1-N1-C2
4	A	1404	XZQ	O7-C32-C33-O8
4	A	1404	XZQ	C33-O8-P1-O10
4	A	1404	XZQ	C34-O14-P2-O12
3	A	1401	XZU	O1-C1-N1-C2
4	A	1404	XZQ	O5-C30-C31-C32
4	A	1404	XZQ	O5-C30-C31-O6
3	A	1401	XZU	C11-C12-C13-C14
3	A	1401	XZU	C20-C12-C13-C14
3	A	1403	XZU	C3-C2-N1-C1
3	A	1403	XZU	C7-C2-N1-C1
4	A	1404	XZQ	P2-O11-P1-O8
4	A	1404	XZQ	C29-C30-C31-C32
4	A	1404	XZQ	P1-O11-P2-O12
4	A	1404	XZQ	P1-O11-P2-O13
4	A	1404	XZQ	C29-C30-C31-O6
4	A	1404	XZQ	C31-C32-C33-O8
4	A	1404	XZQ	O14-C34-C35-O15

There are no ring outliers.

3 monomers are involved in 7 short contacts:

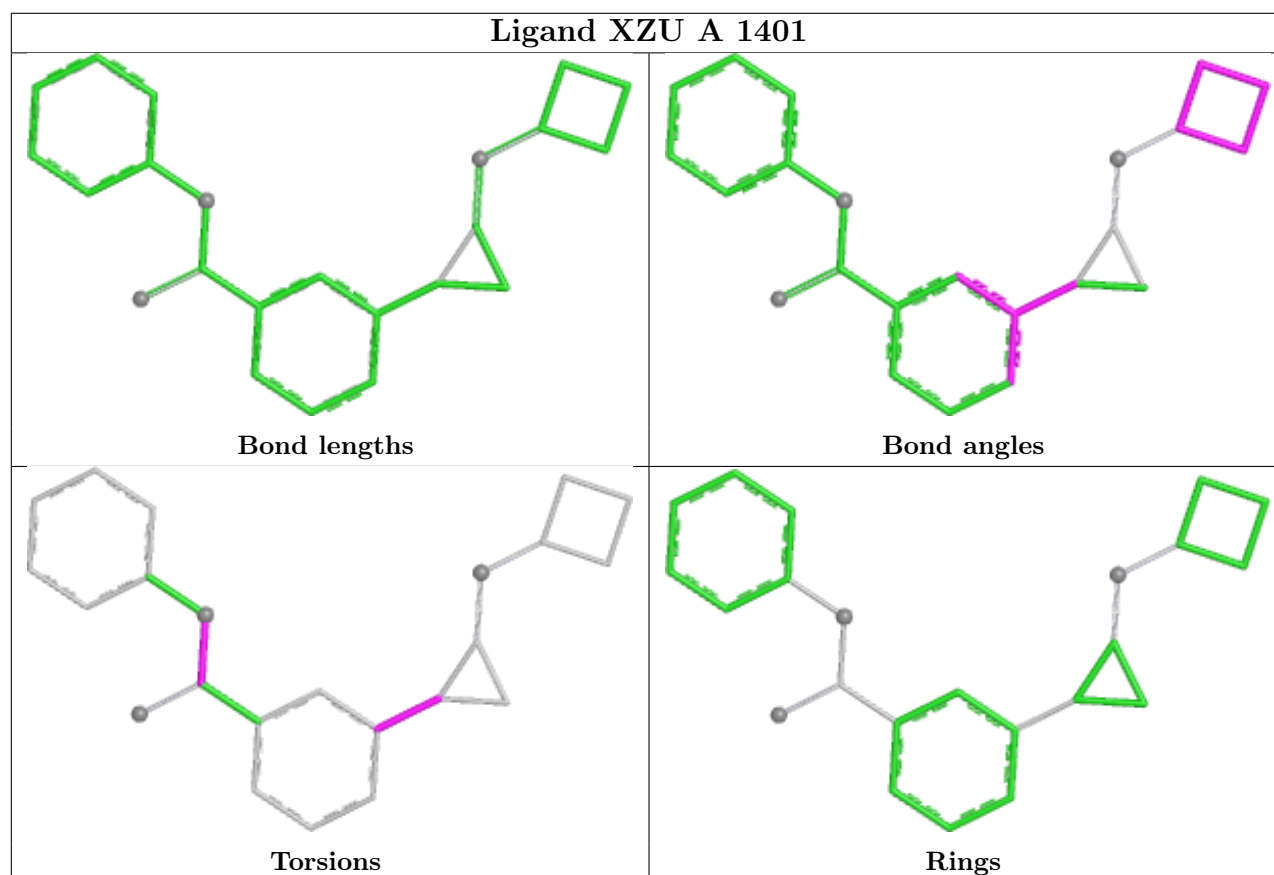
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1401	XZU	1	0

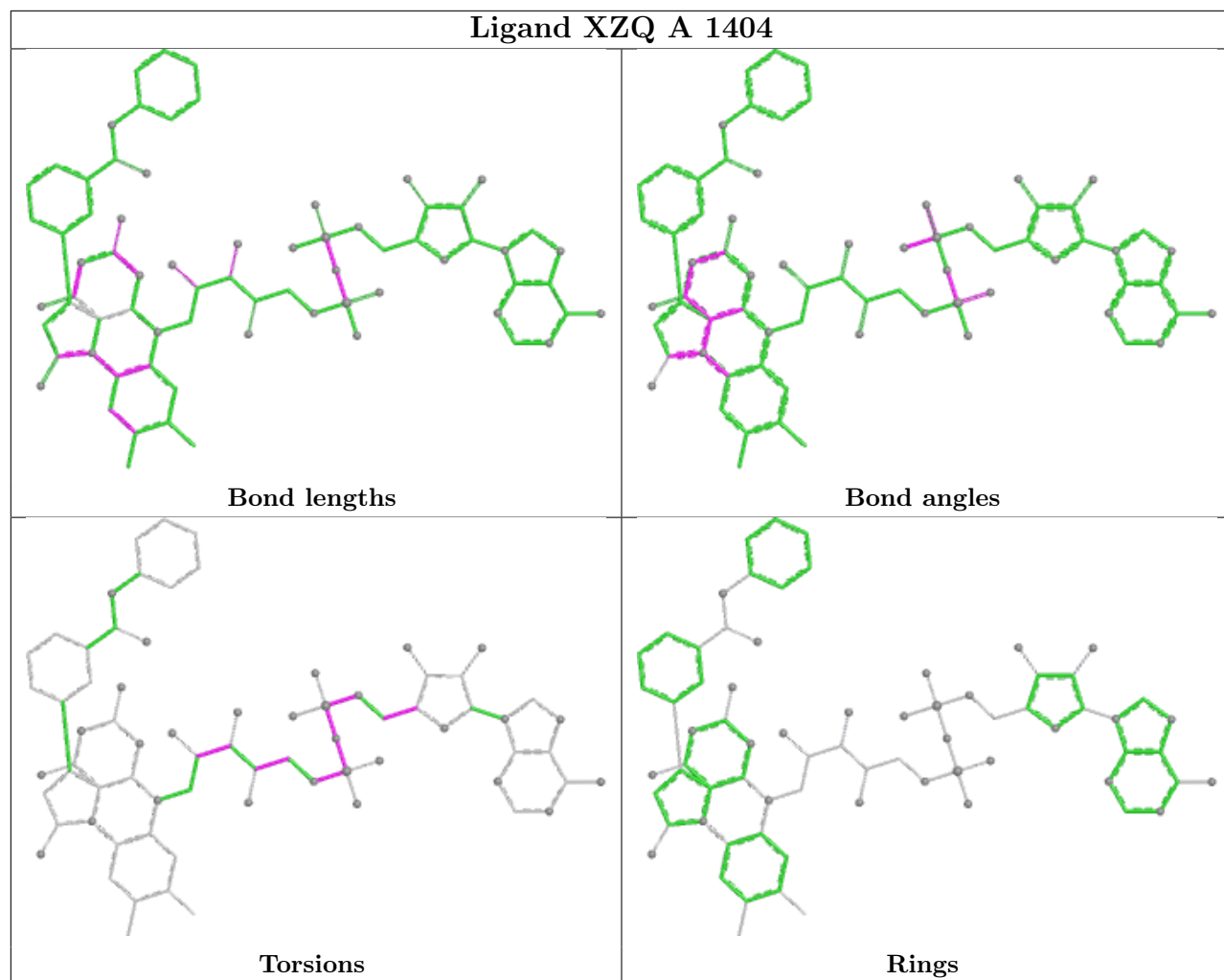
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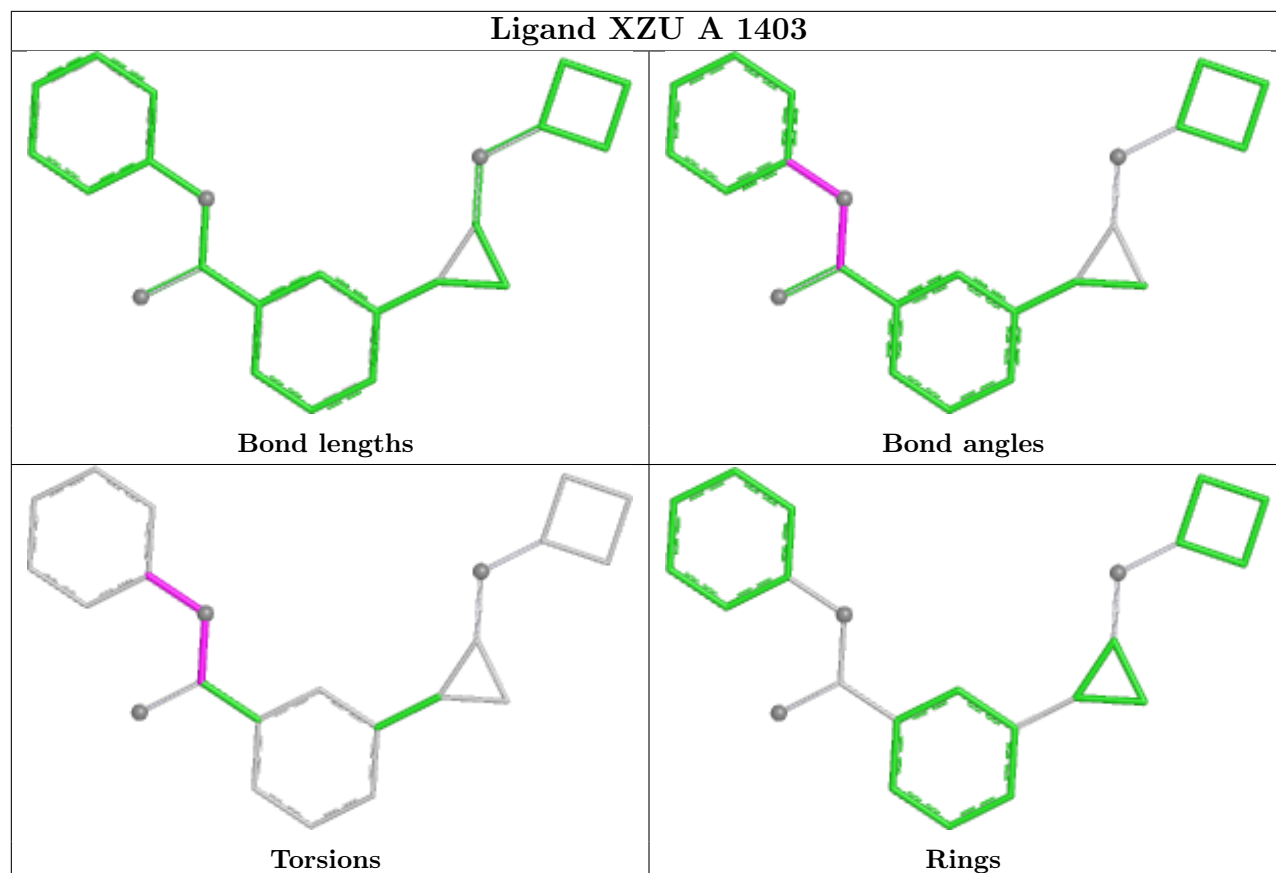
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1404	XZQ	5	0
3	A	1403	XZU	1	0

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.

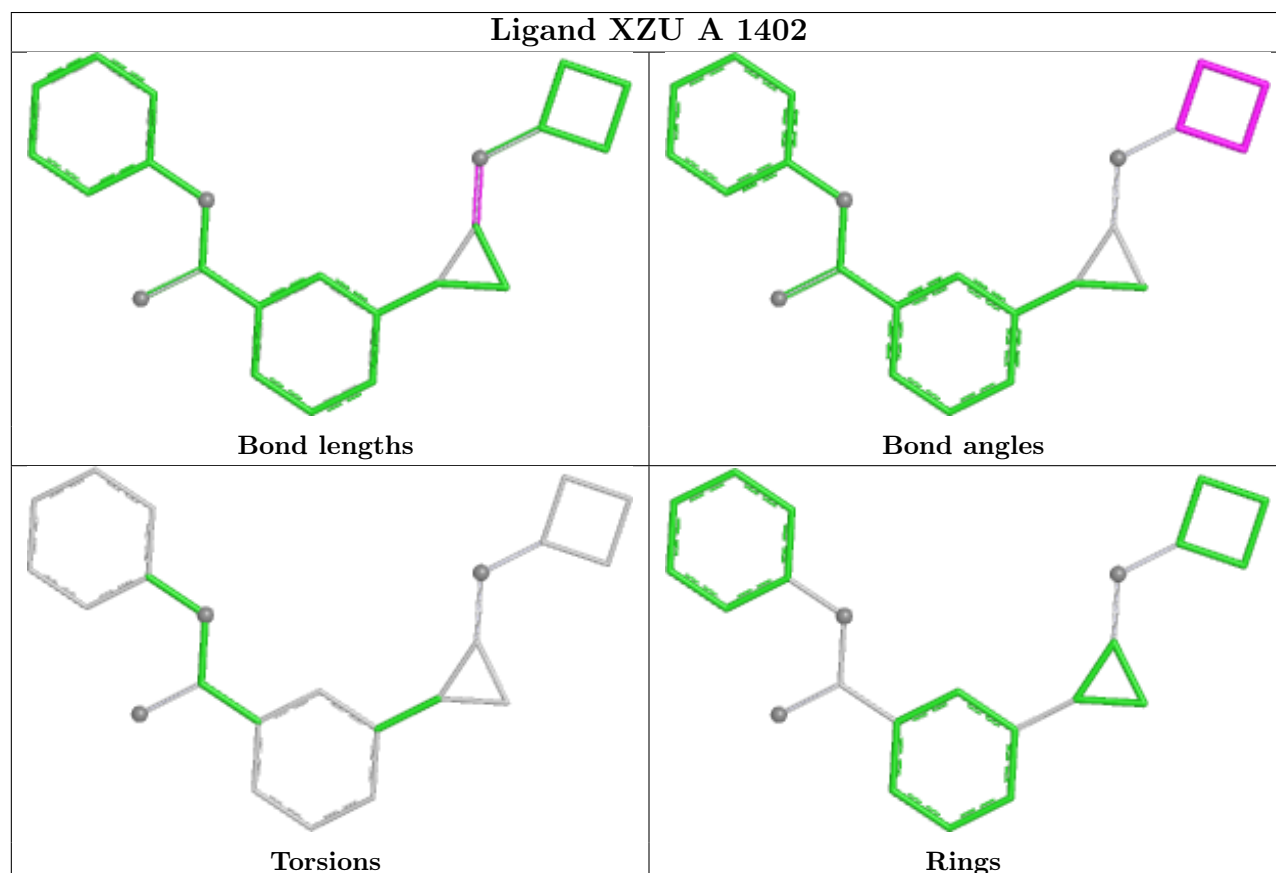




Ligand XZU A 1403



Ligand XZU A 1402



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	666/871 (76%)	0.66	37 (5%)	30 30	58, 84, 117, 135	0
2	B	133/144 (92%)	1.12	20 (15%)	5 5	83, 116, 138, 148	0
All	All	799/1015 (78%)	0.73	57 (7%)	22 22	58, 90, 128, 148	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	836	LEU	7.7
2	B	414	VAL	4.9
1	A	331	ALA	4.7
2	B	376	ILE	4.6
2	B	440	GLU	4.5
1	A	835	THR	4.1
1	A	171	PRO	3.8
1	A	833	MET	3.8
2	B	375	VAL	3.5
1	A	668	ARG	3.4
1	A	358	GLN	3.4
2	B	347	ARG	3.4
1	A	333	VAL	3.3
1	A	453	GLU	3.1
2	B	402	PHE	3.0
2	B	312	LYS	3.0
2	B	439	ALA	3.0
1	A	273	LEU	2.9
1	A	503	LYS	2.9
2	B	340	MET	2.9
1	A	287	GLY	2.9
1	A	507	LYS	2.7
1	A	232	GLU	2.7
1	A	272	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	511	LEU	2.6
2	B	415	VAL	2.6
2	B	378	LYS	2.6
1	A	456	LYS	2.6
1	A	761	TYR	2.4
1	A	740	VAL	2.4
2	B	337	GLN	2.4
1	A	539	ALA	2.4
1	A	275	THR	2.4
1	A	788	GLY	2.4
2	B	308	ARG	2.4
1	A	506	GLU	2.4
1	A	774	ASP	2.3
1	A	791	GLN	2.3
1	A	363	TYR	2.3
1	A	485	ARG	2.3
2	B	422	VAL	2.3
1	A	508	LEU	2.3
2	B	326	ALA	2.2
2	B	413	SER	2.2
1	A	834	TYR	2.2
1	A	778	GLN	2.2
1	A	611	SER	2.2
2	B	425	ARG	2.1
1	A	332	MET	2.1
1	A	373	GLU	2.1
1	A	335	THR	2.1
1	A	271	LYS	2.1
2	B	424	TYR	2.1
1	A	430	HIS	2.1
2	B	371	ARG	2.0
1	A	467	GLU	2.0
2	B	383	TRP	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](#)

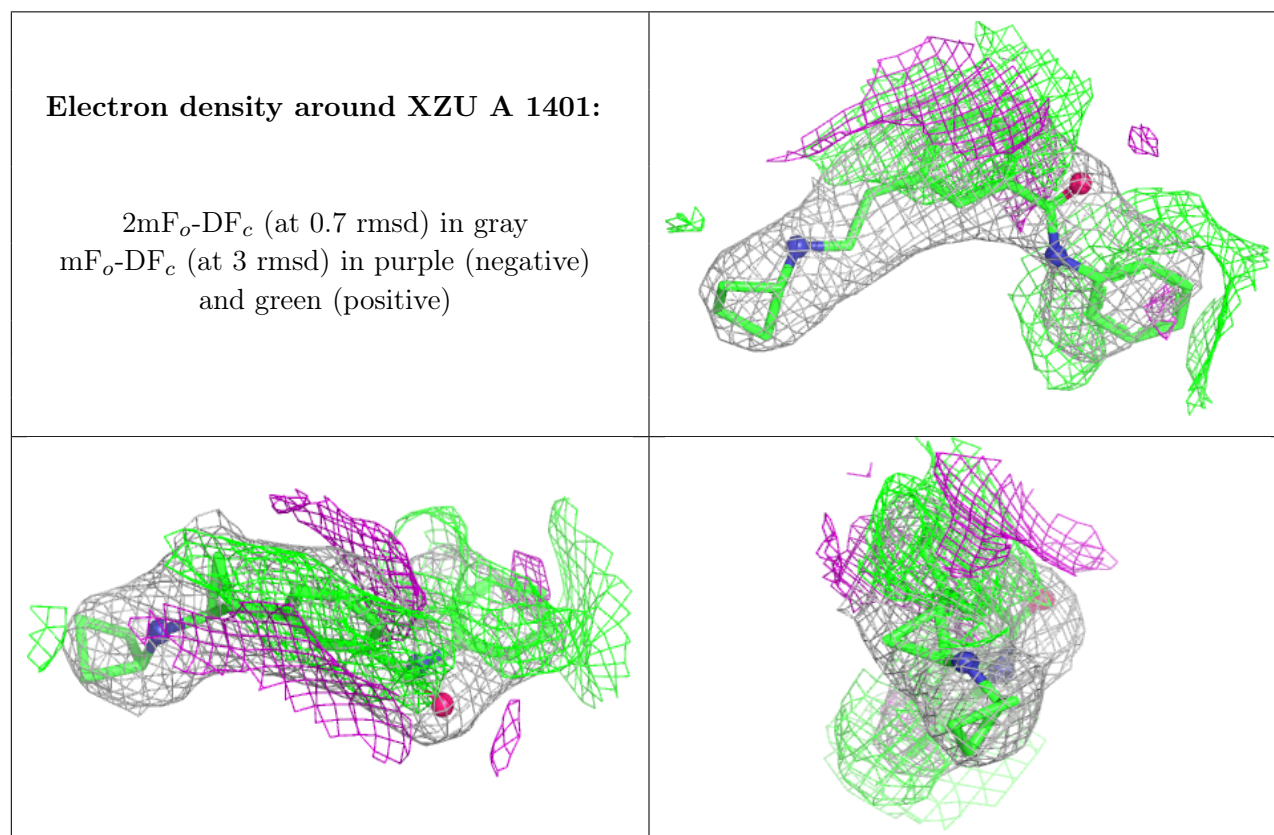
There are no oligosaccharides in this entry.

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.

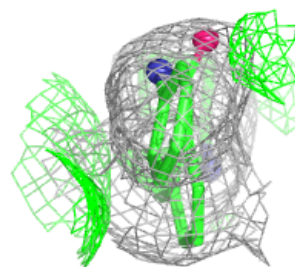
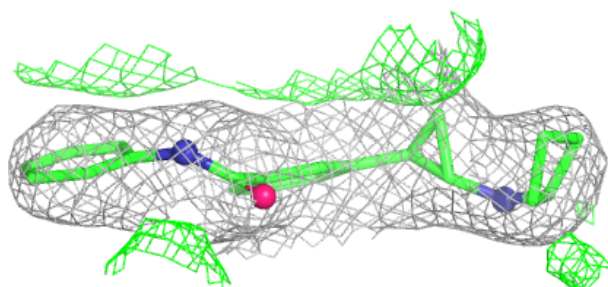
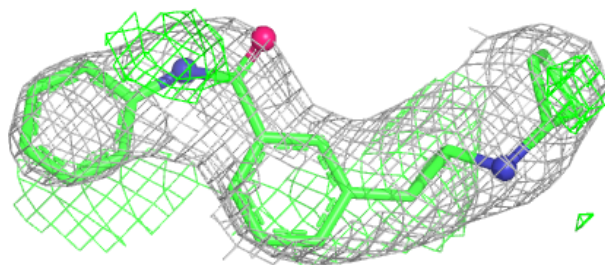
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	XZU	A	1401	23/23	0.87	0.26	70,83,91,98	0
3	XZU	A	1403	23/23	0.87	0.22	88,95,101,108	0
4	XZQ	A	1404	72/72	0.91	0.12	0,0,0,0	0
3	XZU	A	1402	23/23	0.94	0.16	77,83,92,97	0

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.

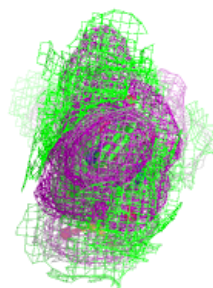
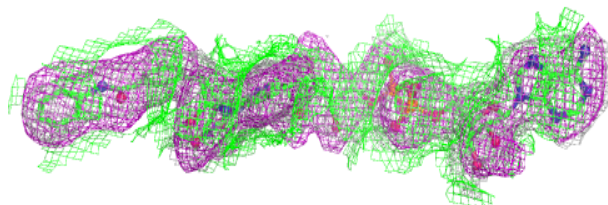
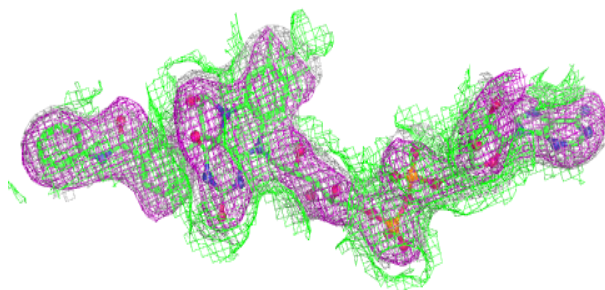


Electron density around XZU A 1403:

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

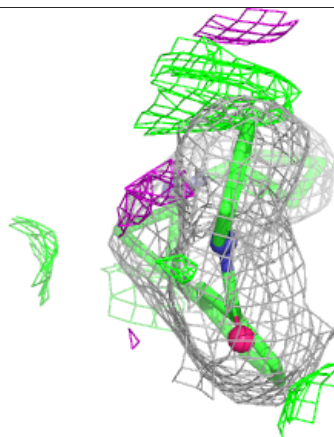
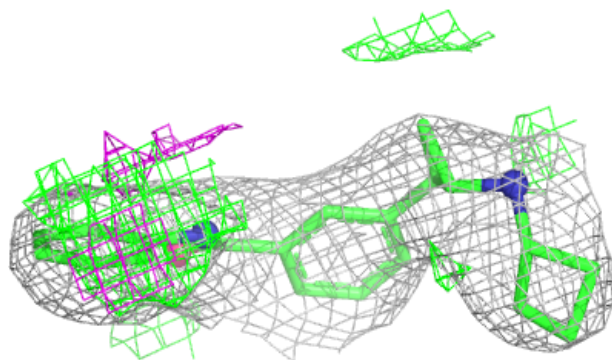
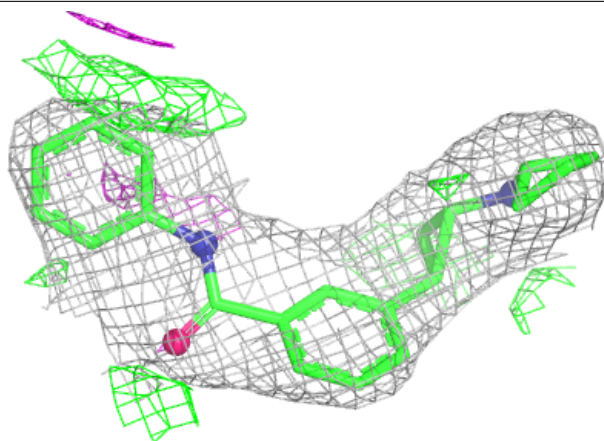
**Electron density around XZQ A 1404:**

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



Electron density around XZU A 1402:

$2mF_o - DF_c$ (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.