Full wwPDB X-ray Structure Validation Report (i)

Nov 2, 2026 – 10:22 AM EST

PDB ID	:	12LU / pdb_000012lu
Title	:	LSD1-CoREST in complex with GSK-LSD1 and peptide PRSALV
Authors	:	Caroli, J.; Mattevi, A.
Deposited on	:	2026-04-12
Resolution	:	3.20 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 (i) 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 (i)) were used in the production of this report:

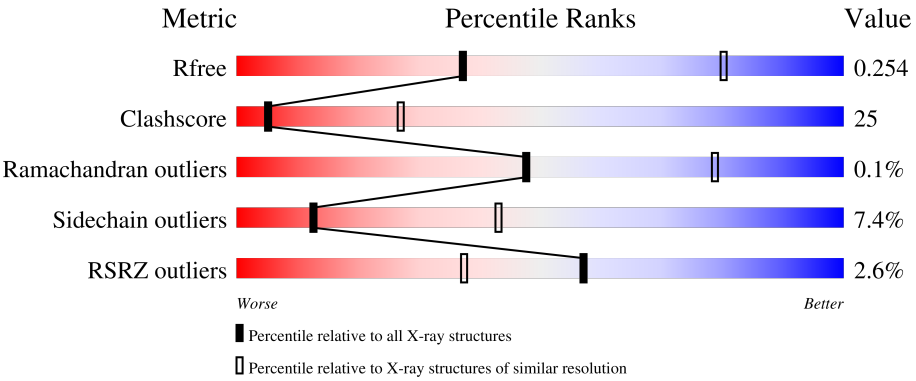
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

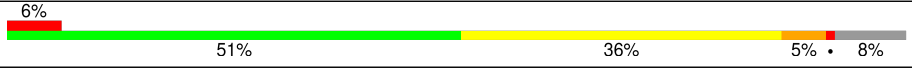
The reported resolution of this entry is 3.20 Å.

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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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 <=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	B	144	
3	E	6	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6400 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			

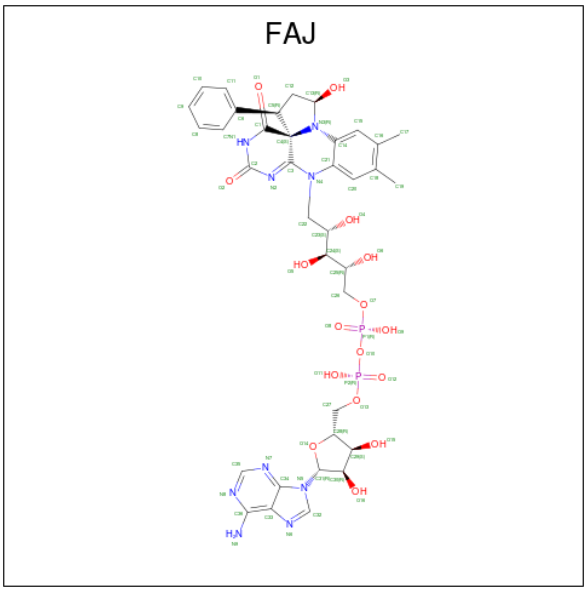
There are 8 discrepancies between the modelled and reference sequences:

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 PRSAL.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	6	Total	C	N	O	0	0	0
			44	28	9	7			

- Molecule 4 is FAD-trans-2-Phenylcyclopropylamine Adduct (CCD ID: FAJ) (formula: $C_{36}H_{43}N_9O_{16}P_2$) (labeled as "Ligand of Interest" by depositor).

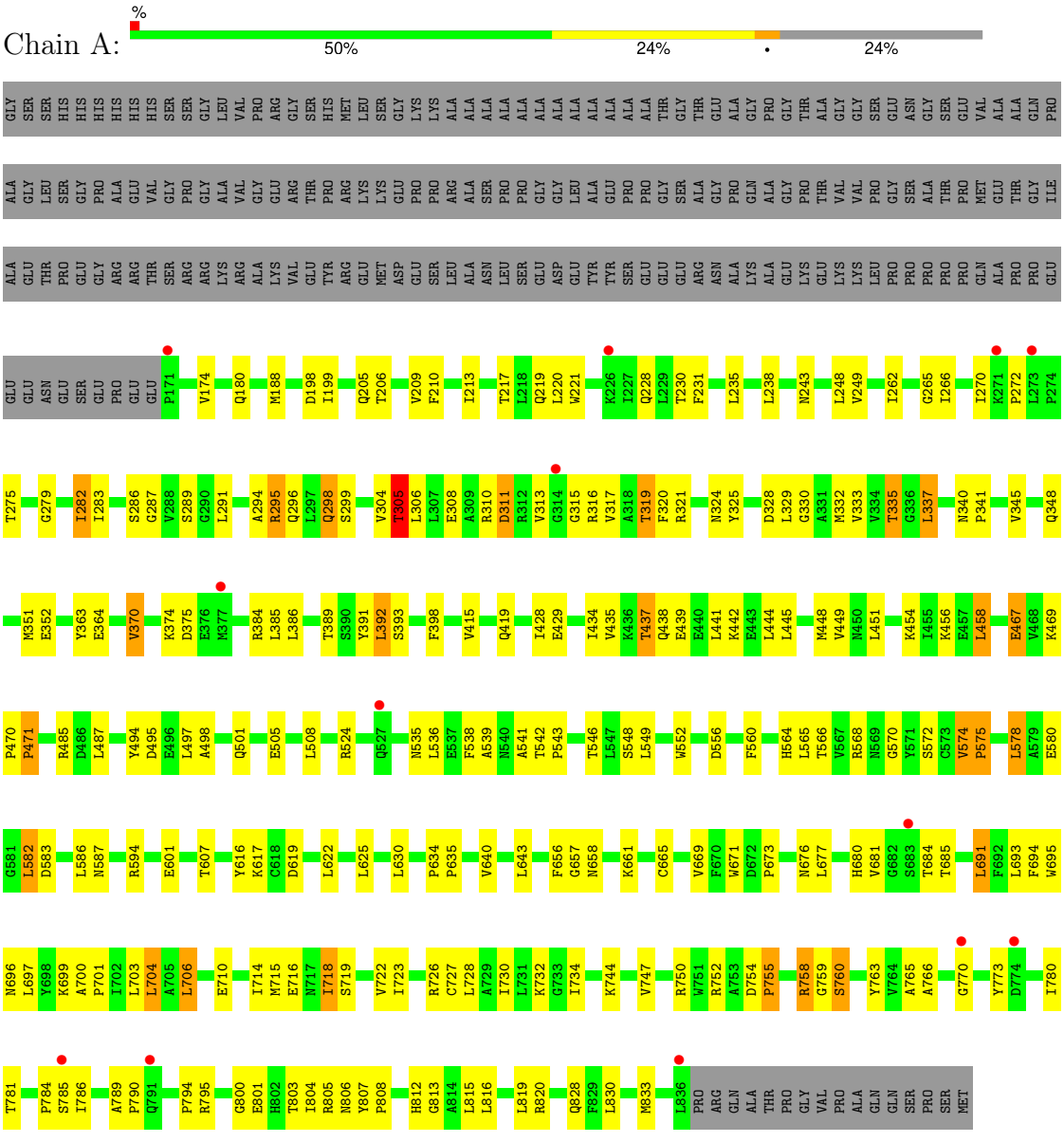


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			63	36	9	16	2		

3 Residue-property plots (i)

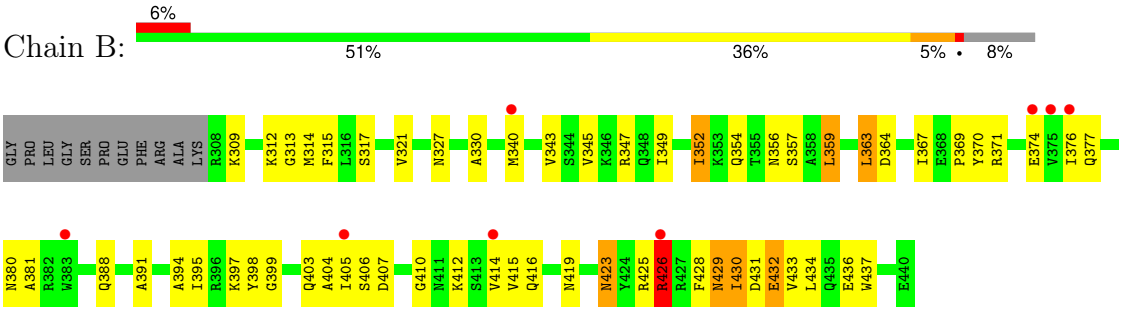
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

Chain B:



• Molecule 3: PRSAL

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	118.28Å 181.01Å 233.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.12 – 3.20 49.12 – 3.20	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.12-3.20) 98.7 (49.12-3.20)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.06 (at 3.19Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.227 , 0.258 0.235 , 0.254	Depositor DCC
R_{free} test set	2000 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å^2)	84.2	Xtriage
Anisotropy	0.624	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 58.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6400	wwPDB-VP
Average B, all atoms (Å^2)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: FAJ

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	0.88	2/5331 (0.0%)	1.07	8/7232 (0.1%)
2	B	0.71	0/1091	1.00	1/1471 (0.1%)
3	E	0.97	0/44	1.62	1/58 (1.7%)
All	All	0.85	2/6466 (0.0%)	1.07	10/8761 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	575	PRO	C-O	-5.77	1.17	1.24
1	A	755	PRO	C-O	-5.53	1.17	1.24

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	315	GLY	CA-C-N	7.48	132.57	120.60
1	A	315	GLY	C-N-CA	7.48	132.57	120.60
1	A	574	VAL	CB-CA-C	-5.92	106.24	114.00
1	A	444	LEU	CA-C-O	-5.58	115.14	121.00
2	B	426	ARG	N-CA-C	-5.52	106.59	113.55
3	E	1	PRO	CA-N-CD	-5.34	104.52	112.00
1	A	279	GLY	CA-C-O	-5.32	117.73	121.88
1	A	454	LYS	N-CA-C	-5.28	105.20	111.69
1	A	305	THR	CB-CA-C	5.12	118.61	109.65
1	A	471	PRO	N-CA-CB	-5.06	97.03	102.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	272	1
2	B	1076	0	1091	73	1
3	E	44	0	52	27	0
4	A	63	0	41	6	0
All	All	6400	0	6436	315	1

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 25.

All (315) 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:325:TYR:CE2	1:A:665:CYS:HB3	1.48	1.49
1:A:535:ASN:HD21	3:E:6:VAL:CG1	1.28	1.45
1:A:539:ALA:CB	3:E:4:ALA:HB3	1.60	1.31
1:A:535:ASN:ND2	3:E:6:VAL:CG1	1.94	1.30
1:A:755:PRO:HA	1:A:758:ARG:NH1	1.56	1.20
1:A:437:THR:CG2	1:A:508:LEU:HD21	1.74	1.18
1:A:535:ASN:HD21	3:E:6:VAL:HG13	1.05	1.17
1:A:438:GLN:OE1	1:A:508:LEU:HD13	1.46	1.15
1:A:556:ASP:OD2	3:E:2:ARG:NH1	1.77	1.15
1:A:539:ALA:HB1	3:E:4:ALA:HB3	1.21	1.13
1:A:691:LEU:CD2	1:A:727:CYS:SG	2.40	1.09
1:A:755:PRO:HA	1:A:758:ARG:HH12	0.96	1.09
1:A:535:ASN:HD21	3:E:6:VAL:HG12	1.09	1.09
1:A:320:PHE:CD1	1:A:747:VAL:HG21	1.88	1.08
1:A:539:ALA:HA	3:E:4:ALA:CB	1.84	1.08
1:A:437:THR:HG21	1:A:508:LEU:CD2	1.85	1.07
1:A:437:THR:HG21	1:A:508:LEU:CG	1.84	1.06
2:B:419:ASN:O	2:B:423:ASN:OD1	1.73	1.05
1:A:325:TYR:HE2	1:A:665:CYS:CB	1.71	1.03
1:A:320:PHE:CE1	1:A:747:VAL:HG21	1.94	1.02
1:A:325:TYR:CE2	1:A:665:CYS:CB	2.41	1.02
1:A:296:GLN:O	1:A:299:SER:HB3	1.60	1.01
1:A:539:ALA:CA	3:E:4:ALA:HB3	1.91	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:535:ASN:ND2	3:E:6:VAL:HG12	1.66	0.99
1:A:754:ASP:OD1	1:A:755:PRO:HD2	1.62	0.99
1:A:691:LEU:HD21	1:A:727:CYS:SG	2.02	0.99
1:A:437:THR:HG22	1:A:508:LEU:HD11	1.44	0.98
1:A:535:ASN:ND2	3:E:6:VAL:HG13	1.67	0.98
1:A:325:TYR:HD2	1:A:665:CYS:SG	1.88	0.97
1:A:695:TRP:HE1	1:A:706:LEU:HD13	1.29	0.96
1:A:320:PHE:CE1	1:A:747:VAL:CG2	2.51	0.94
1:A:291:LEU:CD2	1:A:578:LEU:O	2.15	0.94
1:A:437:THR:CG2	1:A:508:LEU:CD2	2.43	0.94
1:A:437:THR:HG21	1:A:508:LEU:HG	1.48	0.93
1:A:758:ARG:HG2	1:A:758:ARG:HH11	1.32	0.93
1:A:291:LEU:HD22	1:A:578:LEU:O	1.70	0.92
1:A:325:TYR:HD2	1:A:665:CYS:HG	0.99	0.92
1:A:539:ALA:HA	3:E:4:ALA:HB1	1.52	0.91
1:A:469:LYS:HE2	1:A:469:LYS:HA	1.53	0.90
1:A:325:TYR:CD2	1:A:665:CYS:SG	2.63	0.90
1:A:437:THR:CG2	1:A:508:LEU:CG	2.50	0.89
1:A:539:ALA:CA	3:E:4:ALA:CB	2.49	0.88
1:A:437:THR:CG2	1:A:508:LEU:HD11	2.04	0.86
1:A:755:PRO:CA	1:A:758:ARG:HH12	1.85	0.86
1:A:535:ASN:CG	3:E:6:VAL:HG12	2.00	0.86
1:A:392:LEU:HD23	1:A:398:PHE:CD2	2.11	0.85
1:A:656:PHE:CE2	1:A:759:GLY:HA3	2.13	0.83
1:A:716:GLU:OE2	1:A:750:ARG:HA	1.76	0.83
1:A:437:THR:HG22	1:A:508:LEU:HD21	1.59	0.82
1:A:691:LEU:HD22	1:A:727:CYS:SG	2.18	0.82
1:A:539:ALA:CB	3:E:4:ALA:CB	2.53	0.81
1:A:449:VAL:CG2	2:B:363:LEU:HD21	2.08	0.81
1:A:325:TYR:CD2	1:A:665:CYS:HB3	2.16	0.81
1:A:695:TRP:HE1	1:A:706:LEU:CD1	1.92	0.81
1:A:286:SER:OG	1:A:313:VAL:HG12	1.81	0.80
1:A:538:PHE:CE1	1:A:706:LEU:HD21	2.15	0.80
1:A:341:PRO:HG3	1:A:816:LEU:HD21	1.64	0.79
1:A:716:GLU:OE2	1:A:750:ARG:CA	2.30	0.78
1:A:789:ALA:HB1	1:A:790:PRO:HD2	1.66	0.78
1:A:442:LYS:HB2	2:B:356:ASN:OD1	1.83	0.78
1:A:449:VAL:HG22	2:B:363:LEU:HD21	1.64	0.77
1:A:319:THR:HB	1:A:572:SER:HB3	1.66	0.77
1:A:438:GLN:OE1	1:A:508:LEU:CD1	2.31	0.76
1:A:217:THR:HA	1:A:220:LEU:HD12	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:469:LYS:HA	1:A:469:LYS:CE	2.16	0.75
1:A:291:LEU:HD23	1:A:578:LEU:O	1.86	0.75
2:B:399:GLY:N	2:B:437:TRP:NE1	2.34	0.75
1:A:439:GLU:HG2	2:B:352:ILE:CD1	2.18	0.74
2:B:406:SER:OG	2:B:412:LYS:O	2.02	0.74
1:A:188:MET:HG2	1:A:210:PHE:HE2	1.51	0.74
1:A:437:THR:CG2	1:A:508:LEU:CD1	2.65	0.73
1:A:456:LYS:HA	2:B:370:TYR:CE1	2.23	0.73
1:A:458:LEU:HB3	1:A:487:LEU:HD13	1.70	0.73
1:A:535:ASN:OD1	3:E:6:VAL:HG12	1.89	0.73
1:A:535:ASN:CG	3:E:6:VAL:CG1	2.59	0.73
1:A:695:TRP:NE1	1:A:706:LEU:HD13	2.03	0.72
1:A:287:GLY:O	1:A:291:LEU:HD12	1.89	0.72
1:A:754:ASP:OD1	1:A:755:PRO:CD	2.37	0.72
1:A:700:ALA:HB1	1:A:701:PRO:HD2	1.71	0.71
1:A:758:ARG:HH11	1:A:758:ARG:CG	2.03	0.71
1:A:437:THR:HG22	1:A:508:LEU:CD1	2.18	0.70
1:A:619:ASP:O	1:A:795:ARG:NH1	2.23	0.70
2:B:403:GLN:OE1	2:B:414:VAL:HG21	1.92	0.70
1:A:815:LEU:C	1:A:815:LEU:HD23	2.16	0.69
1:A:449:VAL:HG22	2:B:363:LEU:CD2	2.22	0.69
1:A:695:TRP:NE1	1:A:706:LEU:CD1	2.55	0.69
1:A:439:GLU:HG2	2:B:352:ILE:HD13	1.74	0.68
1:A:794:PRO:HD2	1:A:828:GLN:NE2	2.09	0.68
1:A:325:TYR:CD2	1:A:665:CYS:CB	2.77	0.68
1:A:538:PHE:CD1	1:A:706:LEU:CD2	2.76	0.68
1:A:305:THR:OG1	1:A:583:ASP:OD2	2.10	0.68
1:A:539:ALA:O	3:E:1:PRO:HD2	1.94	0.67
1:A:230:THR:HG23	1:A:270:ILE:HD12	1.77	0.67
1:A:495:ASP:OD1	2:B:371:ARG:NH2	2.28	0.66
2:B:415:VAL:O	2:B:419:ASN:ND2	2.24	0.65
1:A:656:PHE:CD2	1:A:759:GLY:HA3	2.31	0.65
1:A:384:ARG:NH2	2:B:312:LYS:O	2.28	0.65
1:A:441:LEU:HD23	2:B:356:ASN:HD22	1.60	0.65
1:A:392:LEU:CD2	1:A:398:PHE:CD2	2.80	0.64
1:A:601:GLU:HB3	1:A:617:LYS:HD3	1.79	0.64
1:A:325:TYR:HE2	1:A:665:CYS:HB3	0.76	0.64
1:A:384:ARG:HH22	2:B:313:GLY:HA3	1.62	0.64
1:A:437:THR:HG21	1:A:508:LEU:HD21	1.48	0.63
1:A:337:LEU:HD23	1:A:337:LEU:N	2.12	0.63
1:A:456:LYS:HA	2:B:370:TYR:HE1	1.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:ILE:HD11	1:A:578:LEU:HD23	1.81	0.63
1:A:287:GLY:C	1:A:291:LEU:HD12	2.24	0.63
2:B:340:MET:HA	2:B:343:VAL:HG12	1.80	0.62
1:A:435:VAL:HG12	2:B:349:ILE:HG12	1.81	0.62
1:A:800:GLY:O	1:A:803:THR:OG1	2.18	0.62
1:A:718:ILE:HG22	1:A:723:ILE:HG13	1.82	0.61
1:A:199:ILE:HG23	1:A:206:THR:HG22	1.81	0.61
1:A:316:ARG:NH1	1:A:656:PHE:HZ	1.97	0.61
1:A:538:PHE:CE1	1:A:706:LEU:CD2	2.84	0.61
2:B:380:ASN:CG	2:B:381:ALA:H	2.09	0.60
1:A:265:GLY:O	1:A:295:ARG:HD2	2.02	0.60
1:A:449:VAL:HG23	2:B:363:LEU:HD21	1.82	0.59
1:A:485:ARG:HG3	2:B:407:ASP:HB2	1.85	0.59
1:A:320:PHE:HB2	1:A:329:LEU:HD11	1.84	0.59
1:A:719:SER:OG	1:A:722:VAL:HG23	2.02	0.59
1:A:437:THR:CG2	1:A:508:LEU:HG	2.24	0.59
2:B:394:ALA:HB1	2:B:405:ILE:HG12	1.85	0.59
2:B:403:GLN:OE1	2:B:403:GLN:HA	2.03	0.59
1:A:536:LEU:HA	3:E:5:LEU:HD11	1.84	0.59
1:A:676:ASN:O	1:A:696:ASN:N	2.30	0.58
1:A:209:VAL:O	1:A:213:ILE:HG13	2.03	0.58
1:A:188:MET:HG2	1:A:210:PHE:CE2	2.37	0.58
1:A:437:THR:CB	1:A:508:LEU:HD21	2.32	0.58
1:A:306:LEU:N	1:A:306:LEU:HD12	2.18	0.58
1:A:794:PRO:HD2	1:A:828:GLN:HE21	1.68	0.58
2:B:399:GLY:H	2:B:437:TRP:CD1	2.22	0.58
1:A:298:GLN:HG3	1:A:298:GLN:O	2.04	0.58
4:A:901:FAJ:H8	3:E:1:PRO:HG3	1.85	0.58
1:A:438:GLN:HB3	2:B:352:ILE:CG2	2.34	0.58
1:A:801:GLU:HG2	4:A:901:FAJ:O5	2.03	0.58
1:A:316:ARG:NH1	1:A:656:PHE:CZ	2.72	0.57
2:B:403:GLN:OE1	2:B:414:VAL:CG2	2.52	0.57
1:A:566:THR:HG21	1:A:697:LEU:HD22	1.86	0.57
1:A:286:SER:O	1:A:291:LEU:HD11	2.04	0.57
1:A:308:GLU:HB3	1:A:586:LEU:HD23	1.87	0.57
1:A:794:PRO:CD	1:A:828:GLN:NE2	2.68	0.56
1:A:317:VAL:HG12	1:A:317:VAL:O	2.04	0.56
2:B:327:ASN:ND2	2:B:330:ALA:HB2	2.20	0.56
1:A:180:GLN:O	1:A:180:GLN:HG3	2.05	0.56
1:A:656:PHE:CE2	1:A:759:GLY:CA	2.89	0.56
1:A:364:GLU:HA	1:A:681:VAL:HB	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:VAL:CG1	2:B:349:ILE:CG1	2.84	0.56
1:A:266:ILE:HD11	1:A:578:LEU:CD2	2.35	0.55
1:A:370:VAL:HG12	1:A:375:ASP:HB2	1.89	0.55
1:A:451:LEU:HD23	1:A:494:TYR:HB2	1.87	0.55
1:A:661:LYS:HD3	1:A:704:LEU:HD21	1.87	0.55
1:A:435:VAL:HG12	2:B:349:ILE:CG1	2.37	0.55
1:A:467:GLU:OE2	1:A:467:GLU:HA	2.07	0.54
1:A:296:GLN:O	1:A:299:SER:CB	2.46	0.54
1:A:728:LEU:O	1:A:732:LYS:HG3	2.08	0.54
2:B:380:ASN:OD1	2:B:381:ALA:N	2.35	0.54
1:A:536:LEU:HA	3:E:5:LEU:CD1	2.36	0.54
1:A:329:LEU:HD21	1:A:747:VAL:CG1	2.38	0.54
1:A:538:PHE:CD1	1:A:706:LEU:HD23	2.42	0.54
1:A:435:VAL:CG1	2:B:349:ILE:HG13	2.38	0.53
2:B:363:LEU:HD23	2:B:363:LEU:N	2.23	0.53
2:B:429:ASN:HD22	2:B:432:GLU:HG3	1.74	0.53
1:A:296:GLN:C	1:A:299:SER:HB3	2.30	0.53
1:A:448:MET:HE1	1:A:498:ALA:HB2	1.91	0.53
2:B:425:ARG:HA	2:B:430:ILE:HG13	1.91	0.53
1:A:316:ARG:O	1:A:330:GLY:HA3	2.09	0.53
1:A:320:PHE:C	1:A:320:PHE:CD2	2.85	0.53
1:A:437:THR:HG22	1:A:508:LEU:CD2	2.28	0.52
1:A:543:PRO:HB3	1:A:710:GLU:HG2	1.91	0.52
1:A:552:TRP:CH2	3:E:2:ARG:HG3	2.45	0.52
4:A:901:FAJ:H8	3:E:1:PRO:CG	2.40	0.52
1:A:174:VAL:HG12	1:A:219:GLN:OE1	2.10	0.52
2:B:425:ARG:HG3	2:B:425:ARG:O	2.10	0.51
1:A:319:THR:HG21	1:A:321:ARG:HE	1.75	0.51
1:A:329:LEU:HD21	1:A:747:VAL:HG11	1.92	0.51
1:A:308:GLU:HA	1:A:308:GLU:OE1	2.10	0.51
1:A:320:PHE:HE1	1:A:747:VAL:CG2	2.19	0.51
1:A:384:ARG:NH2	2:B:313:GLY:HA3	2.25	0.51
1:A:716:GLU:HG2	1:A:750:ARG:HG2	1.93	0.51
1:A:445:LEU:HB2	2:B:359:LEU:HD12	1.92	0.51
1:A:238:LEU:HB2	1:A:243:ASN:HB3	1.92	0.50
2:B:399:GLY:CA	2:B:437:TRP:CE2	2.94	0.50
2:B:388:GLN:HB3	2:B:428:PHE:HE2	1.77	0.50
1:A:661:LYS:HD3	1:A:704:LEU:CD2	2.41	0.50
1:A:319:THR:CG2	1:A:321:ARG:HE	2.24	0.50
1:A:716:GLU:OE2	1:A:750:ARG:CB	2.59	0.50
1:A:805:ARG:O	1:A:808:PRO:HD3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:308:GLU:CB	1:A:586:LEU:HD23	2.41	0.49
1:A:449:VAL:CG2	2:B:363:LEU:CD2	2.86	0.49
1:A:294:ALA:HA	1:A:304:VAL:HG11	1.95	0.49
2:B:399:GLY:HA3	2:B:437:TRP:CE2	2.48	0.49
1:A:296:GLN:HG3	1:A:819:LEU:CD2	2.42	0.49
1:A:501:GLN:O	1:A:505:GLU:HB2	2.13	0.49
1:A:658:ASN:HA	1:A:760:SER:OG	2.12	0.49
1:A:770:GLY:O	1:A:773:TYR:HB2	2.12	0.49
2:B:359:LEU:HD23	2:B:359:LEU:N	2.27	0.49
1:A:296:GLN:HA	1:A:299:SER:HB3	1.93	0.49
1:A:694:PHE:HA	1:A:704:LEU:O	2.12	0.49
2:B:399:GLY:N	2:B:437:TRP:CD1	2.81	0.49
1:A:498:ALA:HA	1:A:501:GLN:HB3	1.94	0.48
2:B:374:GLU:O	2:B:376:ILE:HG13	2.13	0.48
2:B:434:LEU:C	2:B:436:GLU:N	2.70	0.48
1:A:363:TYR:CD2	1:A:734:ILE:HG23	2.48	0.48
1:A:332:MET:HE2	1:A:661:LYS:HZ1	1.79	0.48
1:A:574:VAL:HB	1:A:575:PRO:CD	2.43	0.48
1:A:763:TYR:HE1	1:A:765:ALA:HA	1.78	0.48
1:A:419:GLN:HE22	2:B:315:PHE:H	1.60	0.48
2:B:354:GLN:O	2:B:357:SER:N	2.45	0.48
1:A:286:SER:OG	1:A:286:SER:O	2.22	0.48
2:B:434:LEU:O	2:B:437:TRP:N	2.46	0.48
1:A:578:LEU:HD23	1:A:578:LEU:N	2.29	0.48
1:A:384:ARG:NH1	2:B:312:LYS:O	2.47	0.48
1:A:574:VAL:HB	1:A:575:PRO:HD3	1.95	0.48
1:A:438:GLN:CB	2:B:352:ILE:HG21	2.44	0.47
1:A:456:LYS:HA	2:B:370:TYR:CD1	2.49	0.47
3:E:5:LEU:HD12	3:E:5:LEU:HA	1.68	0.47
1:A:310:ARG:HG3	4:A:901:FAJ:O16	2.14	0.47
1:A:594:ARG:HA	1:A:640:VAL:O	2.14	0.47
2:B:388:GLN:HB3	2:B:428:PHE:CE2	2.49	0.47
1:A:643:LEU:HD23	1:A:643:LEU:HA	1.65	0.47
1:A:684:THR:HG22	1:A:685:THR:N	2.30	0.47
1:A:784:PRO:O	1:A:784:PRO:HG2	2.14	0.47
1:A:235:LEU:HD13	1:A:249:VAL:HG11	1.97	0.47
1:A:758:ARG:NH1	1:A:758:ARG:HG2	2.11	0.47
1:A:270:ILE:O	1:A:272:PRO:HD3	2.15	0.47
1:A:438:GLN:HB3	2:B:352:ILE:HG21	1.96	0.47
1:A:458:LEU:HB3	1:A:487:LEU:CD1	2.42	0.47
1:A:548:SER:O	1:A:552:TRP:HB3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:317:SER:O	2:B:321:VAL:HG23	2.15	0.46
1:A:437:THR:HB	1:A:508:LEU:HD21	1.95	0.46
1:A:803:THR:HB	1:A:820:ARG:HH22	1.80	0.46
2:B:395:ILE:HG22	2:B:433:VAL:HG12	1.96	0.46
1:A:221:TRP:CD1	1:A:262:ILE:HA	2.51	0.46
1:A:391:TYR:HE2	2:B:309:LYS:CA	2.29	0.46
1:A:435:VAL:CG1	2:B:349:ILE:HG12	2.42	0.46
1:A:582:LEU:HD23	1:A:582:LEU:N	2.30	0.46
1:A:658:ASN:CA	1:A:760:SER:OG	2.63	0.46
1:A:311:ASP:OD1	1:A:311:ASP:N	2.31	0.46
1:A:392:LEU:HD23	1:A:398:PHE:HD2	1.76	0.46
1:A:286:SER:HB3	1:A:308:GLU:OE1	2.15	0.46
1:A:345:VAL:O	1:A:348:GLN:N	2.50	0.45
1:A:587:ASN:HB3	1:A:607:THR:OG1	2.17	0.45
1:A:680:HIS:CD2	1:A:730:ILE:HG23	2.51	0.45
1:A:781:THR:HA	1:A:794:PRO:HA	1.99	0.45
1:A:601:GLU:OE1	1:A:617:LYS:HE2	2.16	0.45
2:B:395:ILE:HG22	2:B:433:VAL:CG1	2.47	0.45
1:A:231:PHE:HE1	1:A:249:VAL:HG12	1.81	0.45
1:A:442:LYS:N	2:B:356:ASN:HD21	2.14	0.45
1:A:671:TRP:O	1:A:673:PRO:HD3	2.17	0.44
2:B:399:GLY:N	2:B:437:TRP:CE2	2.86	0.44
1:A:364:GLU:OE2	1:A:524:ARG:NH2	2.49	0.44
1:A:352:GLU:HB3	1:A:568:ARG:HB3	2.00	0.44
1:A:601:GLU:HA	1:A:616:TYR:O	2.18	0.44
1:A:715:MET:HE3	1:A:715:MET:HB3	1.87	0.44
1:A:310:ARG:NH1	1:A:754:ASP:OD2	2.39	0.44
1:A:335:THR:O	1:A:335:THR:OG1	2.13	0.44
1:A:669:VAL:HG13	1:A:671:TRP:CE2	2.53	0.44
1:A:677:LEU:HD11	3:E:6:VAL:CG2	2.47	0.44
1:A:333:VAL:HA	1:A:565:LEU:O	2.17	0.43
1:A:815:LEU:C	1:A:815:LEU:CD2	2.88	0.43
2:B:369:PRO:HB2	2:B:370:TYR:CE2	2.53	0.43
1:A:283:ILE:HG12	1:A:622:LEU:HB3	1.99	0.43
1:A:574:VAL:N	1:A:575:PRO:HD2	2.34	0.43
2:B:426:ARG:H	2:B:426:ARG:HG2	1.62	0.43
1:A:320:PHE:HB2	1:A:329:LEU:CD1	2.46	0.43
1:A:289:SER:OG	4:A:901:FAJ:O8	2.27	0.43
1:A:340:ASN:HA	1:A:560:PHE:CE2	2.53	0.43
1:A:393:SER:HB2	1:A:549:LEU:HG	2.01	0.43
1:A:469:LYS:HE2	1:A:470:PRO:HD3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:388:GLN:O	2:B:391:ALA:HB3	2.18	0.43
1:A:716:GLU:OE2	1:A:750:ARG:C	2.61	0.43
1:A:722:VAL:O	1:A:726:ARG:HG3	2.18	0.43
1:A:248:LEU:HD12	1:A:248:LEU:HA	1.77	0.43
1:A:384:ARG:C	2:B:314:MET:HE3	2.44	0.43
1:A:384:ARG:HB3	2:B:314:MET:CE	2.49	0.43
1:A:335:THR:HB	1:A:564:HIS:CD2	2.53	0.43
1:A:671:TRP:CE3	1:A:703:LEU:HD11	2.53	0.43
1:A:352:GLU:OE2	1:A:699:LYS:HE2	2.19	0.42
1:A:548:SER:HB2	1:A:766:ALA:HA	2.01	0.42
1:A:677:LEU:HD11	3:E:6:VAL:HG23	2.00	0.42
2:B:377:GLN:NE2	2:B:410:GLY:HA3	2.34	0.42
1:A:205:GLN:O	1:A:209:VAL:HG23	2.19	0.42
1:A:282:ILE:HD13	1:A:282:ILE:HA	1.83	0.42
1:A:351:MET:HE2	1:A:351:MET:HB3	1.86	0.42
1:A:812:HIS:O	1:A:816:LEU:HG	2.19	0.42
1:A:744:LYS:HD3	1:A:744:LYS:HA	1.68	0.42
2:B:430:ILE:O	2:B:430:ILE:HG22	2.19	0.42
1:A:552:TRP:CZ3	3:E:2:ARG:HG3	2.55	0.42
1:A:625:LEU:HB2	1:A:630:LEU:HD21	2.02	0.42
2:B:381:ALA:HA	2:B:416:GLN:HE22	1.83	0.42
1:A:321:ARG:NH2	1:A:570:GLY:HA3	2.34	0.42
1:A:594:ARG:HG2	1:A:640:VAL:HB	2.02	0.41
4:A:901:FAJ:H18	4:A:901:FAJ:H17	1.83	0.41
1:A:448:MET:HG2	1:A:497:LEU:HB3	2.02	0.41
1:A:715:MET:CE	1:A:723:ILE:HG12	2.49	0.41
1:A:804:ILE:HD13	1:A:813:GLY:O	2.20	0.41
1:A:542:THR:OG1	1:A:546:THR:OG1	2.32	0.41
1:A:198:ASP:OD1	1:A:198:ASP:N	2.54	0.41
1:A:806:ASN:C	1:A:807:TYR:CD2	2.99	0.41
1:A:715:MET:HE3	1:A:723:ILE:HG12	2.02	0.41
1:A:386:LEU:O	1:A:389:THR:OG1	2.27	0.41
1:A:693:LEU:HD12	1:A:694:PHE:H	1.85	0.41
1:A:228:GLN:HB3	1:A:270:ILE:HG13	2.03	0.41
1:A:568:ARG:HG2	1:A:568:ARG:HH11	1.86	0.41
1:A:180:GLN:O	1:A:180:GLN:CG	2.69	0.40
1:A:321:ARG:HH21	1:A:570:GLY:HA3	1.86	0.40
1:A:541:ALA:O	1:A:657:GLY:HA3	2.21	0.40
1:A:385:LEU:HD23	1:A:415:VAL:HG12	2.02	0.40
1:A:658:ASN:ND2	1:A:752:ARG:HB2	2.36	0.40
2:B:397:LYS:HD3	2:B:398:TYR:CE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:494:TYR:CD2	2:B:367:ILE:HD13	2.57	0.40
1:A:634:PRO:HA	1:A:635:PRO:HD3	1.85	0.40
1:A:485:ARG:HD2	2:B:404:ALA:HB1	2.04	0.40
1:A:758:ARG:NH1	1:A:758:ARG:CG	2.71	0.40
2:B:434:LEU:C	2:B:436:GLU:H	2.29	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:324:ASN:OD1	2:B:354:GLN:OE1[8_455]	2.10	0.10

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%)	639 (96%)	25 (4%)	0	100 100
2	B	131/144 (91%)	114 (87%)	17 (13%)	0	100 100
3	E	4/6 (67%)	3 (75%)	0	1 (25%)	0 0
All	All	799/1021 (78%)	756 (95%)	42 (5%)	1 (0%)	48 79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	5	LEU

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%)	531 (94%)	35 (6%)	16 49
2	B	117/125 (94%)	105 (90%)	12 (10%)	7 28
3	E	5/5 (100%)	1 (20%)	4 (80%)	0 0
All	All	688/845 (81%)	637 (93%)	51 (7%)	13 43

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

Mol	Chain	Res	Type
1	A	275	THR
1	A	282	ILE
1	A	295	ARG
1	A	298	GLN
1	A	305	THR
1	A	311	ASP
1	A	319	THR
1	A	328	ASP
1	A	335	THR
1	A	337	LEU
1	A	370	VAL
1	A	374	LYS
1	A	392	LEU
1	A	428	ILE
1	A	429	GLU
1	A	434	ILE
1	A	437	THR
1	A	458	LEU
1	A	467	GLU
1	A	471	PRO
1	A	578	LEU
1	A	580	GLU
1	A	582	LEU
1	A	691	LEU
1	A	704	LEU
1	A	706	LEU
1	A	714	ILE
1	A	718	ILE
1	A	758	ARG
1	A	760	SER

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Mol	Chain	Res	Type
1	A	780	ILE
1	A	785	SER
1	A	786	ILE
1	A	830	LEU
1	A	833	MET
2	B	345	VAL
2	B	347	ARG
2	B	352	ILE
2	B	359	LEU
2	B	363	LEU
2	B	364	ASP
2	B	423	ASN
2	B	426	ARG
2	B	429	ASN
2	B	430	ILE
2	B	431	ASP
2	B	432	GLU
3	E	2	ARG
3	E	3	SER
3	E	5	LEU
3	E	6	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	296	GLN
1	A	298	GLN
1	A	399	ASN
1	A	461	GLN
1	A	535	ASN
1	A	551	HIS
1	A	633	GLN
1	A	680	HIS
1	A	778	GLN
1	A	828	GLN
2	B	416	GLN
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 (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)

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	FAJ	A	901	-	67,70,70	0.80	2 (2%)	91,109,109	1.32	11 (12%)

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	FAJ	A	901	-	-	5/38/106/106	0/7/8/8

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	FAJ	C13-N3	2.97	1.51	1.46
4	A	901	FAJ	C22-C23	2.18	1.55	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	901	FAJ	C1-C4-C3	-6.12	102.54	113.40
4	A	901	FAJ	C4-N3-C13	-3.66	105.82	109.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	901	FAJ	O7-C26-C25	3.09	117.61	109.36
4	A	901	FAJ	O7-P1-O8	-3.02	96.97	108.94
4	A	901	FAJ	O9-P1-O7	2.91	120.77	107.57
4	A	901	FAJ	O11-P2-O12	2.84	125.66	112.44
4	A	901	FAJ	O9-P1-O10	-2.55	100.37	107.27
4	A	901	FAJ	C14-N3-C13	-2.28	118.69	123.92
4	A	901	FAJ	C22-C23-C24	2.17	115.55	109.66
4	A	901	FAJ	O9-P1-O8	2.15	122.45	112.44
4	A	901	FAJ	C1-N1-C2	-2.06	122.32	125.42

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	901	FAJ	P2-O10-P1-O7
4	A	901	FAJ	C12-C5-C6-C7
4	A	901	FAJ	C23-C24-C25-O6
4	A	901	FAJ	C12-C5-C6-C11
4	A	901	FAJ	P1-O10-P2-O11

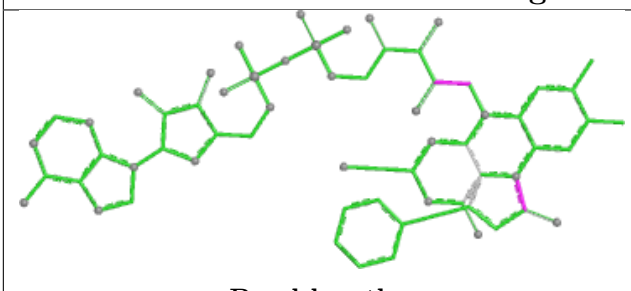
There are no ring outliers.

1 monomer is involved in 6 short contacts:

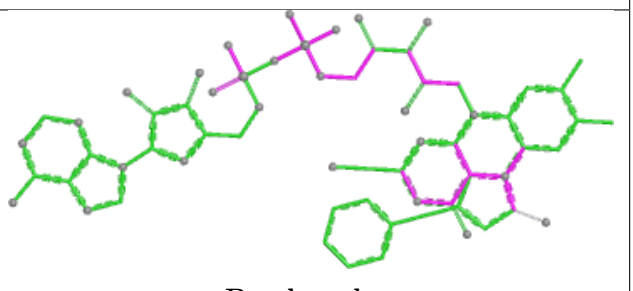
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	901	FAJ	6	0

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 then 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.

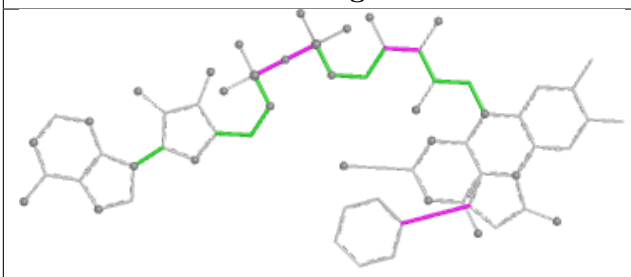
Ligand FAJ A 901



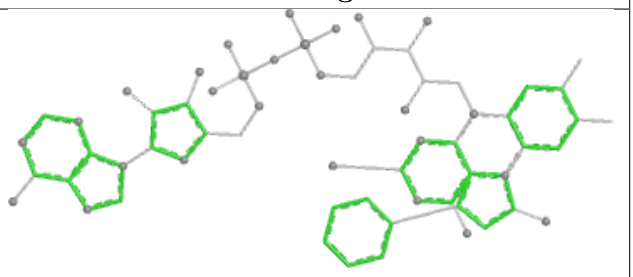
Bond lengths



Bond angles



Torsions



Rings

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, 95^{th} 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%)	0.19	13 (1%) 65 45	47, 79, 111, 134	0
2	B	133/144 (92%)	0.58	8 (6%) 27 17	76, 108, 133, 142	0
3	E	6/6 (100%)	1.07	0 100 100	83, 89, 91, 91	0
All	All	805/1021 (78%)	0.26	21 (2%) 57 37	47, 85, 119, 142	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	836	LEU	5.0
1	A	171	PRO	4.5
1	A	785	SER	3.7
1	A	791	GLN	3.4
2	B	414	VAL	3.0
2	B	426	ARG	3.0
2	B	375	VAL	3.0
1	A	273	LEU	2.9
2	B	405	ILE	2.8
2	B	376	ILE	2.8
1	A	377	MET	2.5
1	A	683	SER	2.4
1	A	314	GLY	2.4
1	A	271	LYS	2.4
2	B	340	MET	2.3
1	A	774	ASP	2.3
2	B	374	GLU	2.3
1	A	527	GLN	2.3
2	B	383	TRP	2.2
1	A	770	GLY	2.0
1	A	226	LYS	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)

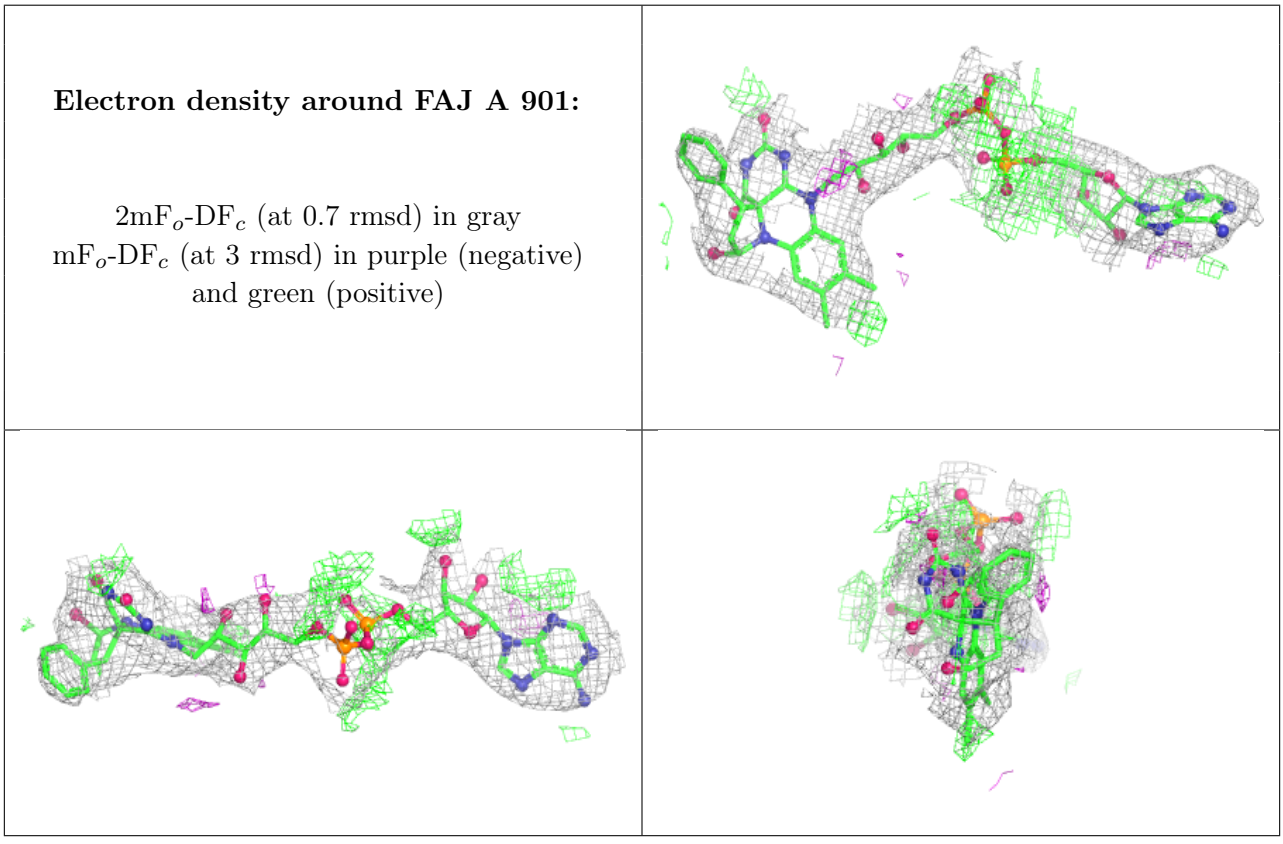
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	FAJ	A	901	63/63	0.96	0.11	51,67,91,97	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 FAJ 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.