



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2024 – 09:56 PM EST

PDB ID : 4MBJ
Title : Human B-Raf Kinase Domain in Complex with an Imidazopyridine-based Inhibitor
Authors : Voegtli, W.C.
Deposited on : 2013-08-19
Resolution : 3.60 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

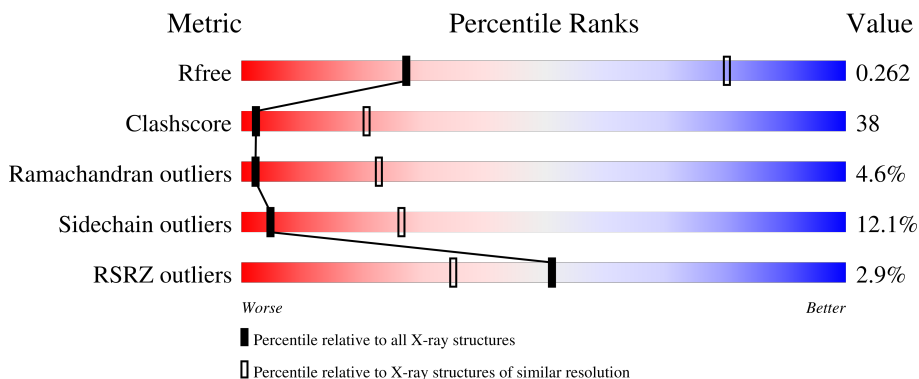
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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}	130704	1257 (3.70-3.50)
Clashscore	141614	1353 (3.70-3.50)
Ramachandran outliers	138981	1307 (3.70-3.50)
Sidechain outliers	138945	1307 (3.70-3.50)
RSRZ outliers	127900	1161 (3.70-3.50)

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	304	
1	B	304	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	DFS	A	801	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4420 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.

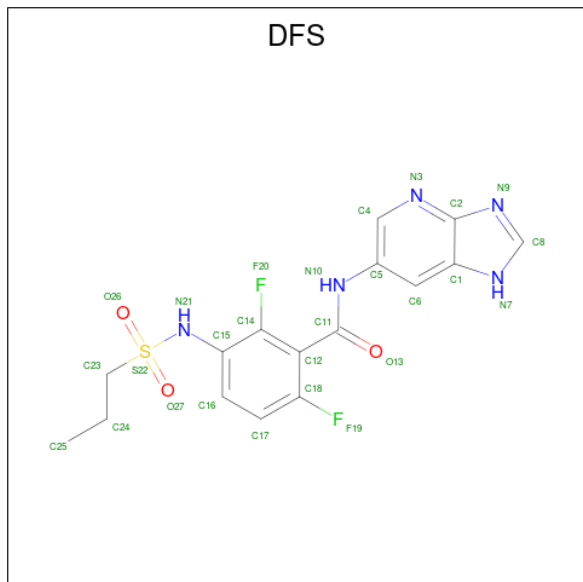
- Molecule 1 is a protein called Serine/threonine-protein kinase B-raf.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	269	Total	C	N	O	S	0	0	0
			2154	1381	376	384	13			
1	B	276	Total	C	N	O	S	0	0	0
			2212	1416	387	396	13			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	420	MET	-	expression tag	UNP P15056
A	421	ASP	-	expression tag	UNP P15056
A	422	ARG	-	expression tag	UNP P15056
A	423	GLY	-	expression tag	UNP P15056
A	424	SER	-	expression tag	UNP P15056
A	425	HIS	-	expression tag	UNP P15056
A	426	HIS	-	expression tag	UNP P15056
A	427	HIS	-	expression tag	UNP P15056
A	428	HIS	-	expression tag	UNP P15056
A	429	HIS	-	expression tag	UNP P15056
A	430	HIS	-	expression tag	UNP P15056
A	431	GLY	-	expression tag	UNP P15056
B	420	MET	-	expression tag	UNP P15056
B	421	ASP	-	expression tag	UNP P15056
B	422	ARG	-	expression tag	UNP P15056
B	423	GLY	-	expression tag	UNP P15056
B	424	SER	-	expression tag	UNP P15056
B	425	HIS	-	expression tag	UNP P15056
B	426	HIS	-	expression tag	UNP P15056
B	427	HIS	-	expression tag	UNP P15056
B	428	HIS	-	expression tag	UNP P15056
B	429	HIS	-	expression tag	UNP P15056
B	430	HIS	-	expression tag	UNP P15056
B	431	GLY	-	expression tag	UNP P15056

- Molecule 2 is 2,6-difluoro-N-(1H-imidazo[4,5-b]pyridin-6-yl)-3-[(propylsulfonyl)amino]benzamide (three-letter code: DFS) (formula: $C_{16}H_{15}F_2N_5O_3S$).

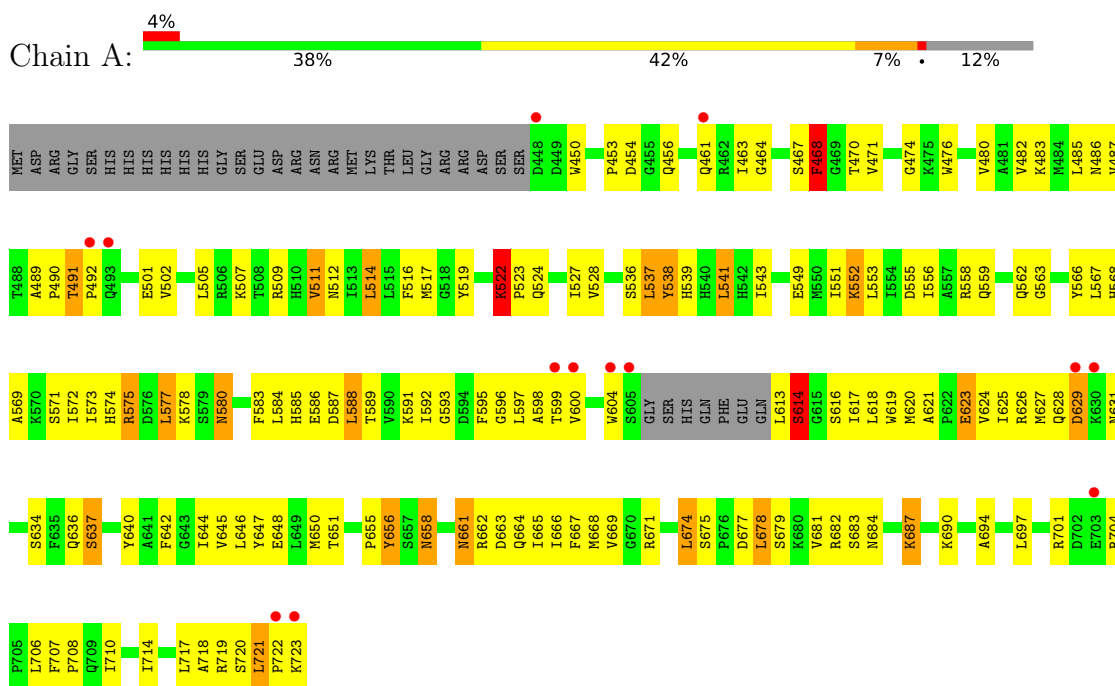


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	F	N	O			S
2	A	1	Total	16	2	5	3	1	0	0
2	B	1	Total	16	2	5	3	1	0	0

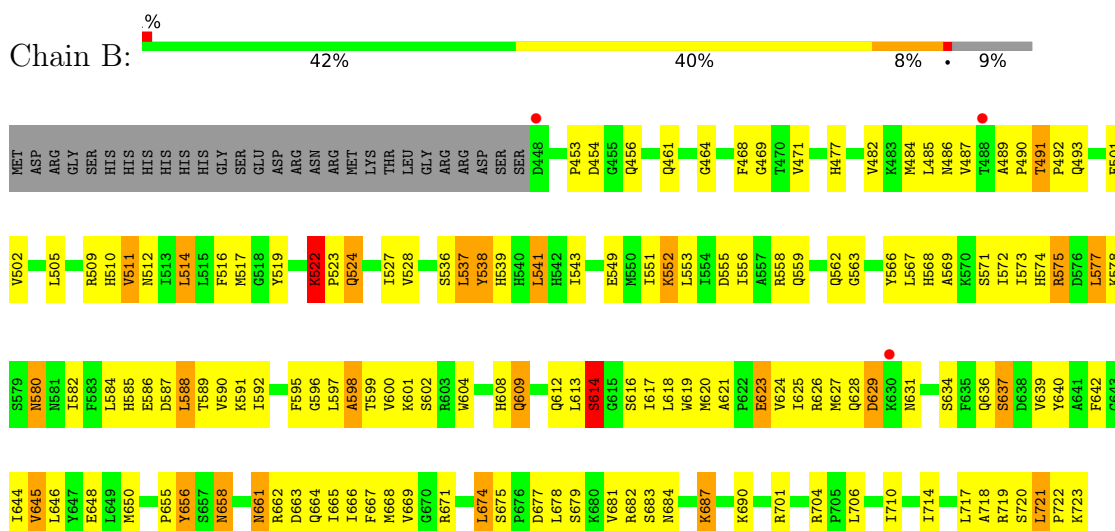
3 Residue-property plots

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: Serine/threonine-protein kinase B-raf



- Molecule 1: Serine/threonine-protein kinase B-raf



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	106.97Å 106.97Å 152.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.60 29.81 – 3.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-3.60) 99.7 (29.81-3.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 3.56Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.276 , 0.328 0.217 , 0.262	Depositor DCC
R_{free} test set	659 reflections (6.15%)	wwPDB-VP
Wilson B-factor (Å ²)	99.3	Xtrriage
Anisotropy	0.067	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4420	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.76% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.46	0/2201	0.59	0/2969
1	B	0.47	0/2262	0.61	0/3052
All	All	0.47	0/4463	0.60	0/6021

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
1	B	0	2
All	All	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	592	ILE	Peptide
1	B	592	ILE	Peptide
1	B	612	GLN	Peptide

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	2154	0	2197	177	0
1	B	2212	0	2244	160	0
2	A	27	0	14	20	0
2	B	27	0	14	7	0
All	All	4420	0	4469	338	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 38.

All (338) 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:580:ASN:HD22	1:A:580:ASN:N	1.42	1.15
1:B:580:ASN:H	1:B:580:ASN:ND2	1.48	1.10
1:B:580:ASN:HD22	1:B:580:ASN:N	1.46	1.06
1:B:454:ASP:HB2	1:B:522:LYS:HD3	1.37	1.05
1:A:454:ASP:HB2	1:A:522:LYS:HD3	1.35	1.03
1:A:580:ASN:H	1:A:580:ASN:ND2	1.44	1.01
1:A:551:ILE:H	1:A:551:ILE:HD12	1.24	0.99
1:B:551:ILE:HD12	1:B:551:ILE:H	1.26	0.97
1:B:658:ASN:HD22	1:B:658:ASN:H	1.10	0.94
1:A:658:ASN:HD22	1:A:658:ASN:H	1.13	0.90
1:B:490:PRO:HA	1:B:491:THR:HB	1.53	0.89
1:A:490:PRO:HA	1:A:491:THR:HB	1.54	0.88
1:A:596:GLY:HA3	2:A:801:DFS:H13	1.53	0.88
1:A:527:ILE:HD12	2:A:801:DFS:H12	1.55	0.87
1:B:490:PRO:HA	1:B:491:THR:CB	2.05	0.86
1:A:490:PRO:HA	1:A:491:THR:CB	2.06	0.85
1:B:505:LEU:HD22	2:B:801:DFS:H10	1.60	0.83
1:A:464:GLY:HA3	1:A:471:VAL:HB	1.61	0.83
1:B:658:ASN:HD22	1:B:658:ASN:N	1.78	0.80
1:A:483:LYS:HZ3	2:A:801:DFS:H8	1.51	0.76
1:A:658:ASN:HD22	1:A:658:ASN:N	1.81	0.76
1:B:501:GLU:HG3	1:B:597:LEU:HA	1.67	0.75
1:A:597:LEU:HG	2:A:801:DFS:H14	1.70	0.74
1:A:509:ARG:HD3	1:B:516:PHE:O	1.87	0.74
1:B:706:LEU:O	1:B:710:ILE:HG13	1.88	0.73
1:B:628:GLN:HG3	1:B:629:ASP:OD2	1.87	0.73
1:B:505:LEU:CD2	2:B:801:DFS:H10	2.18	0.73
1:A:706:LEU:O	1:A:710:ILE:HG13	1.89	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:501:GLU:HG3	1:A:597:LEU:HA	1.70	0.73
1:A:551:ILE:H	1:A:551:ILE:CD1	1.98	0.73
1:A:551:ILE:HD12	1:A:551:ILE:N	2.02	0.72
1:A:650:MET:HE2	1:A:650:MET:HA	1.71	0.72
1:A:628:GLN:HG3	1:A:629:ASP:OD2	1.89	0.71
1:A:617:ILE:HA	1:A:620:MET:HG3	1.73	0.71
1:B:469:GLY:HA2	1:B:484:MET:O	1.91	0.71
2:B:801:DFS:O26	2:B:801:DFS:H7	1.90	0.71
1:A:453:PRO:HD2	1:A:456:GLN:NE2	2.06	0.70
1:A:578:LYS:HD2	1:A:580:ASN:ND2	2.06	0.70
1:A:468:PHE:HE1	1:A:483:LYS:HZ2	1.38	0.68
1:B:551:ILE:HD12	1:B:551:ILE:N	2.05	0.68
1:B:551:ILE:H	1:B:551:ILE:CD1	2.01	0.68
1:B:453:PRO:HD2	1:B:456:GLN:NE2	2.08	0.67
1:A:483:LYS:HD3	2:A:801:DFS:H8	1.76	0.67
1:B:617:ILE:HA	1:B:620:MET:HG3	1.76	0.67
1:A:483:LYS:NZ	2:A:801:DFS:H8	2.10	0.66
1:B:650:MET:HA	1:B:650:MET:HE2	1.78	0.65
2:A:801:DFS:H7	2:A:801:DFS:C23	2.27	0.64
1:B:597:LEU:O	1:B:598:ALA:HB3	1.97	0.64
1:B:600:VAL:O	1:B:600:VAL:HG12	1.97	0.64
1:A:600:VAL:HG12	1:A:600:VAL:O	1.98	0.63
1:A:453:PRO:HD2	1:A:456:GLN:HE21	1.61	0.63
1:B:586:GLU:O	1:B:588:LEU:HD13	1.98	0.63
1:A:483:LYS:CE	2:A:801:DFS:H8	2.29	0.63
1:A:701:ARG:HB2	1:A:704:ARG:NH1	2.14	0.63
1:B:681:VAL:HG21	1:B:690:LYS:HD2	1.81	0.63
1:B:642:PHE:O	1:B:645:VAL:HG23	1.98	0.63
1:B:578:LYS:HD2	1:B:580:ASN:ND2	2.15	0.62
1:B:553:LEU:HB3	1:B:650:MET:HE1	1.81	0.62
1:A:620:MET:HE2	1:A:625:ILE:HG12	1.82	0.62
1:A:483:LYS:HD3	2:A:801:DFS:C17	2.30	0.62
2:A:801:DFS:H7	2:A:801:DFS:H10	1.82	0.61
1:A:597:LEU:O	1:A:598:ALA:HB3	2.00	0.61
2:A:801:DFS:H10	2:A:801:DFS:C16	2.30	0.61
1:A:527:ILE:CD1	2:A:801:DFS:H12	2.29	0.60
1:B:453:PRO:HD2	1:B:456:GLN:HE21	1.64	0.60
1:A:586:GLU:O	1:A:588:LEU:HD13	2.02	0.60
1:B:682:ARG:O	1:B:684:ASN:N	2.34	0.60
1:A:502:VAL:HG11	1:A:519:TYR:CD2	2.37	0.60
1:A:667:PHE:CE1	1:A:671:ARG:HG3	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:501:GLU:HG2	1:A:600:VAL:HG23	1.85	0.59
1:B:617:ILE:HG13	1:B:618:LEU:N	2.16	0.59
2:A:801:DFS:H2	2:A:801:DFS:O13	2.03	0.59
1:A:501:GLU:HB3	1:A:597:LEU:HD23	1.84	0.58
2:B:801:DFS:O26	2:B:801:DFS:C16	2.49	0.58
1:B:620:MET:HE2	1:B:625:ILE:HG12	1.85	0.58
1:A:516:PHE:O	1:B:509:ARG:HD3	2.02	0.58
1:B:501:GLU:HG2	1:B:600:VAL:HG23	1.84	0.58
1:A:483:LYS:CD	2:A:801:DFS:H8	2.33	0.58
1:B:502:VAL:HG11	1:B:519:TYR:CD2	2.39	0.58
1:A:645:VAL:HA	1:A:648:GLU:HG3	1.85	0.58
1:A:490:PRO:CA	1:A:491:THR:HB	2.33	0.57
1:A:617:ILE:HG13	1:A:618:LEU:N	2.18	0.57
1:A:655:PRO:HB2	1:A:656:TYR:CE2	2.40	0.57
1:A:687:LYS:HE3	1:A:687:LYS:HA	1.87	0.57
1:B:556:ILE:HD13	1:B:584:LEU:HD11	1.86	0.57
1:B:664:GLN:O	1:B:668:MET:HG3	2.05	0.57
1:B:710:ILE:O	1:B:714:ILE:HG13	2.05	0.57
1:A:642:PHE:O	1:A:645:VAL:HG23	2.05	0.56
1:A:467:SER:O	1:A:468:PHE:HB3	2.05	0.56
1:A:682:ARG:O	1:A:684:ASN:N	2.38	0.56
1:B:687:LYS:HE3	1:B:687:LYS:HA	1.88	0.56
1:A:574:HIS:CD2	1:A:595:PHE:HB3	2.40	0.56
1:B:613:LEU:O	1:B:614:SER:HB2	2.06	0.56
1:B:684:ASN:H	1:B:684:ASN:ND2	2.04	0.56
1:B:701:ARG:HB2	1:B:704:ARG:NH1	2.21	0.56
1:A:553:LEU:HB3	1:A:650:MET:HE1	1.88	0.55
1:A:647:TYR:O	1:A:651:THR:OG1	2.20	0.55
1:B:634:SER:O	1:B:637:SER:HB3	2.07	0.55
1:A:464:GLY:HA3	1:A:471:VAL:CB	2.35	0.55
1:A:491:THR:N	1:A:492:PRO:HD3	2.22	0.54
1:B:567:LEU:HD22	1:B:572:ILE:HG21	1.90	0.54
1:A:710:ILE:O	1:A:714:ILE:HG13	2.07	0.54
1:B:667:PHE:CE1	1:B:671:ARG:HG3	2.43	0.54
1:B:589:THR:HG21	1:B:591:LYS:HE2	1.89	0.54
1:B:468:PHE:HB2	1:B:608:HIS:CB	2.38	0.54
1:B:512:ASN:ND2	1:B:559:GLN:HB3	2.23	0.54
1:B:580:ASN:ND2	1:B:580:ASN:N	2.20	0.54
1:A:556:ILE:HD13	1:A:584:LEU:HD11	1.89	0.54
1:B:490:PRO:CA	1:B:491:THR:HB	2.32	0.54
1:B:501:GLU:HB3	1:B:597:LEU:HD23	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:621:ALA:HB1	1:A:704:ARG:NH2	2.23	0.53
1:A:568:HIS:O	1:A:571:SER:N	2.42	0.53
1:A:684:ASN:ND2	1:A:684:ASN:H	2.07	0.53
1:B:721:LEU:HB3	1:B:722:PRO:HD3	1.90	0.53
1:B:655:PRO:HB2	1:B:656:TYR:CE2	2.43	0.53
1:B:490:PRO:CA	1:B:491:THR:CB	2.83	0.53
1:B:656:TYR:HE1	1:B:674:LEU:HD22	1.73	0.53
1:B:574:HIS:O	1:B:575:ARG:HB2	2.08	0.53
1:B:623:GLU:OE1	1:B:704:ARG:NH2	2.40	0.53
1:A:681:VAL:HG21	1:A:690:LYS:HD2	1.89	0.53
1:B:577:LEU:HD23	1:B:578:LYS:H	1.74	0.53
1:B:619:TRP:CD1	1:B:644:ILE:HB	2.44	0.53
1:A:514:LEU:HD22	2:A:801:DFS:F20	1.99	0.52
1:A:721:LEU:HB3	1:A:722:PRO:HD3	1.91	0.52
1:B:491:THR:N	1:B:492:PRO:HD3	2.24	0.52
1:B:553:LEU:HD13	1:B:650:MET:HE2	1.90	0.52
1:B:558:ARG:O	1:B:562:GLN:N	2.36	0.52
1:B:454:ASP:CB	1:B:522:LYS:HD3	2.26	0.52
1:B:568:HIS:O	1:B:571:SER:N	2.41	0.52
1:B:636:GLN:NE2	1:B:706:LEU:HD21	2.25	0.52
1:A:553:LEU:HD13	1:A:650:MET:HE2	1.91	0.52
1:B:718:ALA:C	1:B:720:SER:H	2.14	0.52
1:A:589:THR:HG21	1:A:591:LYS:HE2	1.92	0.52
1:A:578:LYS:HD2	1:A:580:ASN:CG	2.30	0.51
1:A:567:LEU:HD22	1:A:572:ILE:HG21	1.93	0.51
1:A:644:ILE:O	1:A:648:GLU:HG3	2.11	0.51
1:A:485:LEU:N	1:A:485:LEU:HD22	2.26	0.51
1:A:580:ASN:N	1:A:580:ASN:ND2	2.17	0.51
1:A:619:TRP:CD1	1:A:644:ILE:HB	2.45	0.51
1:B:645:VAL:HA	1:B:648:GLU:HG3	1.91	0.51
1:A:577:LEU:HD23	1:A:578:LYS:H	1.75	0.51
1:A:623:GLU:OE1	1:A:704:ARG:NH2	2.43	0.51
1:B:501:GLU:OE1	1:B:599:THR:HB	2.10	0.51
1:B:644:ILE:O	1:B:648:GLU:HG3	2.11	0.51
1:A:454:ASP:CB	1:A:522:LYS:HD3	2.25	0.51
1:A:490:PRO:CA	1:A:491:THR:CB	2.83	0.51
1:B:628:GLN:HG3	1:B:629:ASP:N	2.26	0.50
1:A:625:ILE:HG23	1:A:666:ILE:HG23	1.92	0.50
1:A:628:GLN:HG3	1:A:629:ASP:N	2.26	0.50
1:B:658:ASN:H	1:B:658:ASN:ND2	1.91	0.50
1:A:677:ASP:O	1:A:679:SER:N	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:634:SER:O	1:A:637:SER:HB3	2.12	0.50
2:A:801:DFS:C23	2:A:801:DFS:C16	2.90	0.50
1:A:664:GLN:O	1:A:665:ILE:C	2.49	0.50
1:B:658:ASN:N	1:B:658:ASN:ND2	2.51	0.50
1:A:562:GLN:NE2	1:B:477:HIS:C	2.65	0.50
1:B:588:LEU:H	1:B:588:LEU:HD22	1.77	0.50
1:B:620:MET:HE3	1:B:624:VAL:HG12	1.93	0.50
1:B:621:ALA:HB1	1:B:704:ARG:NH2	2.27	0.50
1:A:658:ASN:N	1:A:658:ASN:ND2	2.54	0.49
1:A:577:LEU:O	1:A:578:LYS:HB3	2.11	0.49
1:A:588:LEU:H	1:A:588:LEU:HD22	1.76	0.49
1:A:536:SER:O	1:A:539:HIS:N	2.45	0.49
1:B:626:ARG:HG3	1:B:626:ARG:O	2.13	0.49
1:A:617:ILE:CA	1:A:620:MET:HG3	2.40	0.49
1:A:656:TYR:HE1	1:A:674:LEU:HD22	1.76	0.49
1:B:522:LYS:HB2	1:B:523:PRO:CD	2.42	0.49
1:B:656:TYR:CE1	1:B:674:LEU:HD22	2.47	0.49
1:A:491:THR:O	1:A:491:THR:HG22	2.12	0.49
1:A:536:SER:O	1:A:537:LEU:C	2.50	0.49
1:A:718:ALA:C	1:A:720:SER:H	2.15	0.49
1:A:522:LYS:HB2	1:A:523:PRO:CD	2.42	0.49
1:A:617:ILE:HA	1:A:620:MET:CG	2.42	0.49
1:A:674:LEU:HD23	1:A:675:SER:N	2.27	0.49
1:B:596:GLY:N	2:B:801:DFS:O27	2.43	0.49
1:A:613:LEU:O	1:A:614:SER:HB2	2.12	0.49
1:A:634:SER:H	1:A:637:SER:HB3	1.78	0.49
1:B:536:SER:O	1:B:539:HIS:N	2.45	0.49
1:B:626:ARG:O	1:B:628:GLN:N	2.46	0.49
1:B:674:LEU:HD23	1:B:675:SER:N	2.28	0.49
1:A:559:GLN:O	1:A:562:GLN:HB3	2.13	0.48
1:A:578:LYS:HD2	1:A:580:ASN:HD21	1.77	0.48
1:B:597:LEU:O	1:B:598:ALA:CB	2.59	0.48
1:A:538:TYR:CD1	1:A:538:TYR:C	2.86	0.48
1:A:501:GLU:OE1	1:A:599:THR:HB	2.13	0.48
1:B:577:LEU:O	1:B:578:LYS:HB3	2.13	0.48
1:A:596:GLY:HA3	2:A:801:DFS:C24	2.35	0.48
1:B:621:ALA:HB2	1:B:640:TYR:CB	2.44	0.48
1:A:578:LYS:CE	1:A:616:SER:HB3	2.44	0.47
1:A:658:ASN:H	1:A:658:ASN:ND2	1.96	0.47
1:A:661:ASN:OD1	1:A:661:ASN:C	2.53	0.47
1:B:578:LYS:CE	1:B:616:SER:HB3	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:574:HIS:CD2	1:B:595:PHE:HB3	2.49	0.47
1:A:464:GLY:CA	1:A:471:VAL:HB	2.37	0.47
1:A:553:LEU:HD13	1:A:650:MET:CE	2.44	0.47
1:B:468:PHE:HB2	1:B:608:HIS:HB3	1.97	0.47
1:A:573:ILE:HD13	1:A:634:SER:HA	1.96	0.47
1:A:586:GLU:O	1:A:587:ASP:HB2	2.14	0.47
1:A:626:ARG:O	1:A:626:ARG:HG3	2.15	0.47
1:B:553:LEU:HD13	1:B:650:MET:CE	2.45	0.47
1:A:636:GLN:NE2	1:A:706:LEU:HD21	2.30	0.47
1:A:656:TYR:CE1	1:A:674:LEU:HD22	2.50	0.47
1:B:485:LEU:N	1:B:485:LEU:HD22	2.30	0.47
1:B:623:GLU:H	1:B:623:GLU:HG3	1.33	0.46
1:B:464:GLY:HA3	1:B:471:VAL:HB	1.97	0.46
1:B:491:THR:HG22	1:B:491:THR:O	2.15	0.46
1:B:522:LYS:HB2	1:B:523:PRO:HD3	1.96	0.46
1:B:661:ASN:C	1:B:661:ASN:OD1	2.52	0.46
1:A:569:ALA:CB	1:B:477:HIS:CE1	2.98	0.46
1:A:701:ARG:CB	1:A:704:ARG:NH1	2.78	0.46
1:B:536:SER:O	1:B:537:LEU:C	2.54	0.46
1:B:634:SER:H	1:B:637:SER:HB3	1.80	0.46
1:B:556:ILE:CD1	1:B:584:LEU:HD11	2.46	0.46
1:A:597:LEU:CG	2:A:801:DFS:H14	2.42	0.46
1:A:701:ARG:HB2	1:A:704:ARG:HH12	1.81	0.46
1:B:617:ILE:CA	1:B:620:MET:HG3	2.45	0.46
1:A:665:ILE:O	1:A:669:VAL:HG23	2.16	0.46
1:B:538:TYR:CD1	1:B:538:TYR:C	2.89	0.46
1:B:559:GLN:O	1:B:562:GLN:HB3	2.16	0.46
1:B:625:ILE:HG23	1:B:666:ILE:HG23	1.97	0.45
1:A:620:MET:HE3	1:A:620:MET:HB3	1.65	0.45
1:A:623:GLU:H	1:A:623:GLU:HG3	1.35	0.45
1:B:568:HIS:O	1:B:569:ALA:C	2.55	0.45
1:A:597:LEU:O	1:A:598:ALA:CB	2.62	0.45
1:B:468:PHE:CZ	1:B:485:LEU:HD12	2.51	0.45
1:B:573:ILE:HD13	1:B:634:SER:HA	1.99	0.45
1:B:642:PHE:CE2	1:B:646:LEU:HD11	2.52	0.45
1:B:602:SER:HA	1:B:609:GLN:NE2	2.32	0.45
1:A:505:LEU:CD1	1:A:527:ILE:HD13	2.46	0.44
1:A:522:LYS:HB2	1:A:523:PRO:HD3	1.98	0.44
1:B:492:PRO:HA	1:B:493:GLN:HA	1.60	0.44
1:A:537:LEU:CD1	1:A:541:LEU:HD22	2.47	0.44
1:A:661:ASN:OD1	1:A:664:GLN:HG2	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:664:GLN:O	1:A:668:MET:HG3	2.17	0.44
1:A:668:MET:CB	1:A:674:LEU:HB2	2.47	0.44
1:B:589:THR:CG2	1:B:591:LYS:HE2	2.47	0.44
1:B:639:VAL:O	1:B:642:PHE:HB3	2.16	0.44
1:B:665:ILE:O	1:B:669:VAL:HG23	2.17	0.44
1:A:650:MET:HE2	1:A:650:MET:CA	2.46	0.44
1:B:537:LEU:HD12	1:B:537:LEU:O	2.16	0.44
1:B:677:ASP:O	1:B:679:SER:N	2.51	0.44
1:A:464:GLY:O	1:A:470:THR:HG23	2.17	0.44
1:A:501:GLU:CG	1:A:597:LEU:HA	2.44	0.44
1:B:501:GLU:CG	1:B:597:LEU:HA	2.42	0.44
1:B:668:MET:HB3	1:B:674:LEU:HB2	1.99	0.44
1:A:589:THR:CG2	1:A:591:LYS:HE2	2.47	0.44
1:A:483:LYS:HD3	2:A:801:DFS:C18	2.47	0.44
1:A:511:VAL:O	1:A:511:VAL:CG2	2.65	0.44
1:A:578:LYS:NZ	1:A:616:SER:HB3	2.33	0.44
1:A:626:ARG:O	1:A:628:GLN:N	2.50	0.44
1:B:668:MET:CB	1:B:674:LEU:HB2	2.47	0.44
1:A:555:ASP:OD1	1:A:558:ARG:NH1	2.51	0.44
1:A:621:ALA:HB2	1:A:640:TYR:CB	2.47	0.44
1:B:645:VAL:O	1:B:648:GLU:HB2	2.18	0.44
1:B:505:LEU:CD1	1:B:527:ILE:HD13	2.47	0.43
1:A:556:ILE:CD1	1:A:584:LEU:HD11	2.48	0.43
1:A:568:HIS:O	1:A:569:ALA:C	2.54	0.43
1:B:684:ASN:H	1:B:684:ASN:HD22	1.66	0.43
1:A:586:GLU:OE1	1:A:586:GLU:HA	2.18	0.43
1:A:485:LEU:O	1:A:487:VAL:HG13	2.18	0.43
1:A:620:MET:HE3	1:A:624:VAL:HG12	1.99	0.43
1:A:655:PRO:HB2	1:A:656:TYR:CD2	2.53	0.43
1:A:717:LEU:C	1:A:717:LEU:HD12	2.38	0.43
1:A:717:LEU:HD12	1:A:717:LEU:O	2.18	0.43
1:B:514:LEU:HB3	2:B:801:DFS:H14	2.01	0.43
1:B:580:ASN:H	1:B:580:ASN:HD22	0.65	0.43
1:B:585:HIS:C	1:B:587:ASP:H	2.21	0.43
1:A:644:ILE:HD13	1:A:697:LEU:HD21	1.99	0.43
1:B:512:ASN:HD21	1:B:559:GLN:HB3	1.83	0.43
1:B:505:LEU:HD11	1:B:527:ILE:HD13	2.00	0.43
1:B:514:LEU:HD23	2:B:801:DFS:H12	2.01	0.43
1:B:664:GLN:O	1:B:665:ILE:C	2.56	0.43
1:B:490:PRO:HA	1:B:491:THR:OG1	2.18	0.43
1:B:582:ILE:HG23	1:B:590:VAL:HG13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:619:TRP:HA	1:B:644:ILE:HG13	2.00	0.43
1:B:563:GLY:O	1:B:566:TYR:HB3	2.18	0.43
1:B:636:GLN:HB3	1:B:704:ARG:HH11	1.84	0.43
1:B:655:PRO:HB2	1:B:656:TYR:CD2	2.54	0.43
1:A:552:LYS:O	1:A:555:ASP:N	2.52	0.43
1:B:586:GLU:O	1:B:587:ASP:HB2	2.19	0.43
1:A:674:LEU:CD2	1:A:675:SER:N	2.82	0.43
1:A:512:ASN:ND2	1:A:559:GLN:HB3	2.33	0.42
1:A:600:VAL:HG13	1:A:604:TRP:CE2	2.55	0.42
1:A:617:ILE:HG13	1:A:618:LEU:HD12	2.01	0.42
1:A:642:PHE:CE2	1:A:646:LEU:HD11	2.54	0.42
1:B:681:VAL:HG21	1:B:690:LYS:CD	2.47	0.42
1:A:490:PRO:HA	1:A:491:THR:OG1	2.18	0.42
1:A:491:THR:N	1:A:492:PRO:CD	2.82	0.42
1:A:678:LEU:HD13	1:A:694:ALA:HB2	2.01	0.42
1:B:578:LYS:NZ	1:B:616:SER:HB3	2.34	0.42
1:B:620:MET:HE3	1:B:620:MET:HB3	1.71	0.42
1:A:574:HIS:O	1:A:575:ARG:HB2	2.20	0.42
1:A:661:ASN:OD1	1:A:661:ASN:O	2.37	0.42
1:A:450:TRP:CH2	1:B:509:ARG:HD2	2.54	0.42
1:A:468:PHE:HE1	1:A:483:LYS:NZ	2.12	0.42
1:A:665:ILE:H	1:A:665:ILE:HG12	1.69	0.42
1:A:563:GLY:O	1:A:566:TYR:HB3	2.19	0.42
1:B:600:VAL:HG13	1:B:604:TRP:CE2	2.54	0.42
1:A:467:SER:O	1:A:468:PHE:CB	2.67	0.42
1:A:684:ASN:H	1:A:684:ASN:HD22	1.67	0.42
1:B:614:SER:O	1:B:617:ILE:HG23	2.19	0.42
1:B:620:MET:CE	1:B:625:ILE:HA	2.50	0.42
1:B:510:HIS:ND1	1:B:511:VAL:N	2.68	0.42
1:B:617:ILE:HG13	1:B:618:LEU:HD12	2.02	0.42
1:A:476:TRP:HB3	1:A:480:VAL:HG11	2.02	0.41
1:B:524:GLN:HE21	1:B:524:GLN:HB2	1.53	0.41
1:B:586:GLU:OE1	1:B:586:GLU:HA	2.19	0.41
1:A:558:ARG:O	1:A:562:GLN:N	2.38	0.41
1:B:717:LEU:HD12	1:B:717:LEU:C	2.40	0.41
1:A:553:LEU:HB3	1:A:650:MET:CE	2.49	0.41
1:A:707:PHE:N	1:A:708:PRO:CD	2.83	0.41
1:B:552:LYS:O	1:B:555:ASP:N	2.54	0.41
1:A:668:MET:HB3	1:A:674:LEU:HB2	2.03	0.41
1:B:578:LYS:HD2	1:B:580:ASN:HD21	1.84	0.41
1:A:541:LEU:HD12	1:A:541:LEU:HA	1.81	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:569:ALA:HB2	1:B:477:HIS:HE1	1.86	0.41
1:B:482:VAL:HG22	1:B:528:VAL:HG13	2.03	0.41
1:B:541:LEU:HD12	1:B:541:LEU:HA	1.81	0.41
1:A:467:SER:O	1:A:468:PHE:HD2	2.03	0.41
1:A:543:ILE:HD13	1:A:543:ILE:HA	1.73	0.41
1:A:583:PHE:HE1	1:A:593:GLY:HA2	1.86	0.41
1:B:543:ILE:HA	1:B:543:ILE:HD13	1.70	0.41
1:B:717:LEU:HD12	1:B:717:LEU:O	2.21	0.41
1:B:485:LEU:O	1:B:487:VAL:HG13	2.20	0.41
1:A:636:GLN:HB3	1:A:704:ARG:HH11	1.86	0.40
1:B:552:LYS:O	1:B:555:ASP:HB3	2.21	0.40
1:B:595:PHE:CD2	1:B:595:PHE:C	2.95	0.40
1:A:580:ASN:HD22	1:A:580:ASN:H	0.61	0.40
1:B:555:ASP:OD1	1:B:558:ARG:NH1	2.54	0.40
1:A:474:GLY:N	1:A:480:VAL:O	2.51	0.40
1:A:482:VAL:HG22	1:A:528:VAL:HG13	2.04	0.40
1:A:585:HIS:C	1:A:587:ASP:H	2.24	0.40
1:B:601:LYS:HG3	1:B:608:HIS:ND1	2.36	0.40
1:A:483:LYS:HD3	2:A:801:DFS:F19	2.11	0.40
1:A:625:ILE:HG22	1:A:626:ARG:N	2.35	0.40
1:B:511:VAL:O	1:B:511:VAL:CG2	2.69	0.40
1:A:471:VAL:HG22	1:A:483:LYS:HG3	2.04	0.40
1:B:491:THR:N	1:B:492:PRO:CD	2.84	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	265/304 (87%)	221 (83%)	31 (12%)	13 (5%)	2 21
1	B	274/304 (90%)	224 (82%)	38 (14%)	12 (4%)	2 23

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	539/608 (89%)	445 (83%)	69 (13%)	25 (5%)	2	23

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	486	ASN
1	A	683	SER
1	B	486	ASN
1	B	683	SER
1	A	614	SER
1	A	721	LEU
1	B	614	SER
1	B	721	LEU
1	A	489	ALA
1	A	491	THR
1	A	719	ARG
1	B	489	ALA
1	B	627	MET
1	B	719	ARG
1	B	491	THR
1	B	598	ALA
1	B	678	LEU
1	A	627	MET
1	A	629	ASP
1	A	678	LEU
1	B	629	ASP
1	A	468	PHE
1	A	522	LYS
1	B	522	LYS
1	A	463	ILE

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	237/268 (88%)	208 (88%)	29 (12%)	5	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	243/268 (91%)	214 (88%)	29 (12%)	5	27
All	All	480/536 (90%)	422 (88%)	58 (12%)	5	26

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	461	GLN
1	A	468	PHE
1	A	507	LYS
1	A	511	VAL
1	A	514	LEU
1	A	517	MET
1	A	522	LYS
1	A	524	GLN
1	A	537	LEU
1	A	538	TYR
1	A	541	LEU
1	A	549	GLU
1	A	552	LYS
1	A	575	ARG
1	A	577	LEU
1	A	580	ASN
1	A	588	LEU
1	A	614	SER
1	A	623	GLU
1	A	631	ASN
1	A	637	SER
1	A	656	TYR
1	A	658	ASN
1	A	661	ASN
1	A	662	ARG
1	A	663	ASP
1	A	674	LEU
1	A	687	LYS
1	A	723	LYS
1	B	461	GLN
1	B	511	VAL
1	B	514	LEU
1	B	517	MET
1	B	522	LYS
1	B	524	GLN
1	B	537	LEU

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Mol	Chain	Res	Type
1	B	538	TYR
1	B	541	LEU
1	B	549	GLU
1	B	552	LYS
1	B	575	ARG
1	B	577	LEU
1	B	580	ASN
1	B	588	LEU
1	B	609	GLN
1	B	614	SER
1	B	623	GLU
1	B	631	ASN
1	B	637	SER
1	B	645	VAL
1	B	656	TYR
1	B	658	ASN
1	B	661	ASN
1	B	662	ARG
1	B	663	ASP
1	B	674	LEU
1	B	687	LYS
1	B	723	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	456	GLN
1	A	461	GLN
1	A	496	GLN
1	A	500	ASN
1	A	524	GLN
1	A	562	GLN
1	A	580	ASN
1	A	631	ASN
1	A	658	ASN
1	A	684	ASN
1	B	456	GLN
1	B	477	HIS
1	B	496	GLN
1	B	500	ASN
1	B	524	GLN
1	B	580	ASN

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Mol	Chain	Res	Type
1	B	609	GLN
1	B	631	ASN
1	B	658	ASN
1	B	684	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
2	DFS	B	801	-	27,29,29	3.53	6 (22%)	32,42,42	1.65	8 (25%)
2	DFS	A	801	-	27,29,29	3.44	13 (48%)	32,42,42	1.61	5 (15%)

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
2	DFS	B	801	-	-	3/17/17/17	0/3/3/3
2	DFS	A	801	-	-	7/17/17/17	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	DFS	S22-N21	-16.50	1.29	1.62
2	A	801	DFS	C15-N21	-13.65	1.20	1.42
2	B	801	DFS	C15-N21	-4.69	1.35	1.42
2	A	801	DFS	C15-C14	4.69	1.46	1.38
2	A	801	DFS	S22-N21	4.53	1.71	1.62
2	A	801	DFS	C16-C15	-3.74	1.33	1.39
2	A	801	DFS	O26-S22	3.49	1.48	1.43
2	A	801	DFS	C4-N3	3.22	1.37	1.31
2	A	801	DFS	O27-S22	3.20	1.48	1.43
2	A	801	DFS	C17-C18	3.02	1.44	1.37
2	B	801	DFS	C4-N3	2.88	1.36	1.31
2	B	801	DFS	C2-N3	-2.79	1.33	1.37
2	B	801	DFS	O26-S22	2.77	1.47	1.43
2	A	801	DFS	C2-N9	2.72	1.39	1.34
2	B	801	DFS	O27-S22	2.51	1.47	1.43
2	A	801	DFS	C12-C18	-2.26	1.35	1.39
2	A	801	DFS	C6-C1	-2.20	1.38	1.41
2	A	801	DFS	C12-C14	-2.15	1.35	1.38
2	A	801	DFS	C1-C2	-2.01	1.35	1.40

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	DFS	C18-C12-C14	4.85	119.52	115.84
2	A	801	DFS	C18-C12-C14	4.52	119.26	115.84
2	B	801	DFS	C23-S22-N21	-3.95	101.14	106.77
2	A	801	DFS	C16-C15-N21	-3.56	113.76	120.88
2	A	801	DFS	C23-S22-N21	-2.95	102.55	106.77
2	B	801	DFS	C5-C6-C1	-2.94	117.29	120.50
2	A	801	DFS	C4-N3-C2	-2.71	113.97	116.69
2	B	801	DFS	C4-N3-C2	-2.61	114.07	116.69
2	B	801	DFS	C5-C4-N3	2.12	126.10	124.49
2	A	801	DFS	C16-C15-C14	2.09	120.88	117.80
2	B	801	DFS	C16-C15-C14	2.08	120.86	117.80
2	B	801	DFS	C18-C12-C11	-2.04	119.41	122.27
2	B	801	DFS	C6-C1-C2	-2.02	119.09	121.12

There are no chirality outliers.

All (10) torsion outliers are listed below:

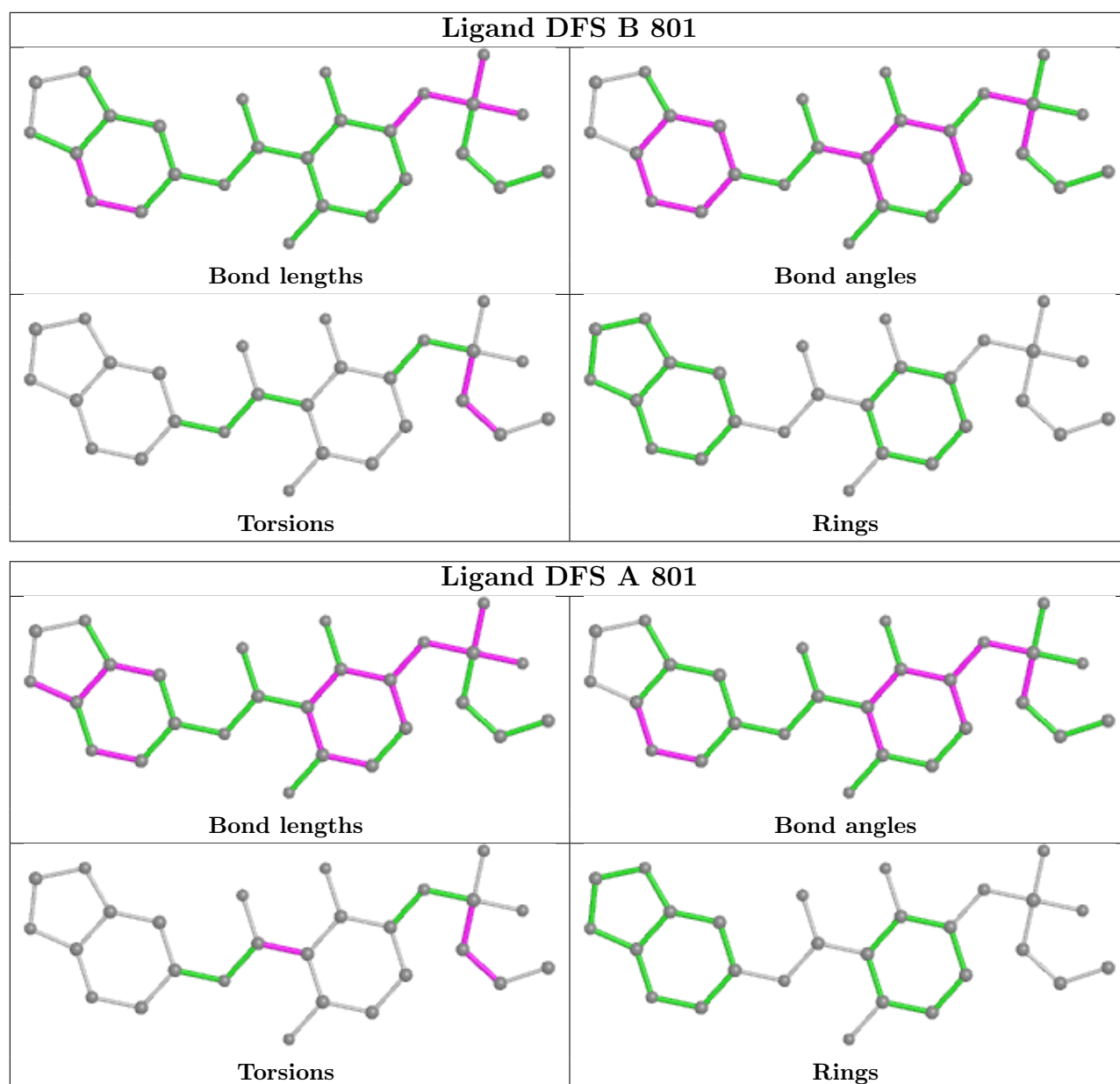
Mol	Chain	Res	Type	Atoms
2	A	801	DFS	N10-C11-C12-C14
2	A	801	DFS	N10-C11-C12-C18
2	A	801	DFS	O13-C11-C12-C14
2	A	801	DFS	O13-C11-C12-C18
2	A	801	DFS	C24-C23-S22-N21
2	A	801	DFS	C24-C23-S22-O27
2	B	801	DFS	C24-C23-S22-N21
2	B	801	DFS	C24-C23-S22-O26
2	B	801	DFS	S22-C23-C24-C25
2	A	801	DFS	S22-C23-C24-C25

There are no ring outliers.

2 monomers are involved in 27 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	DFS	7	0
2	A	801	DFS	20	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	269/304 (88%)	-0.15	13 (4%) 30 19	24, 55, 115, 138	0
1	B	276/304 (90%)	-0.37	3 (1%) 80 68	22, 55, 115, 138	0
All	All	545/608 (89%)	-0.26	16 (2%) 51 35	22, 55, 115, 138	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	723	LYS	3.5
1	A	722	PRO	3.4
1	A	492	PRO	2.9
1	A	448	ASP	2.8
1	B	630	LYS	2.8
1	A	605	SER	2.7
1	A	630	LYS	2.7
1	A	604	TRP	2.5
1	A	600	VAL	2.5
1	A	599	THR	2.4
1	B	488	THR	2.4
1	A	493	GLN	2.4
1	A	461	GLN	2.3
1	A	629	ASP	2.1
1	A	703	GLU	2.1
1	B	448	ASP	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](#)

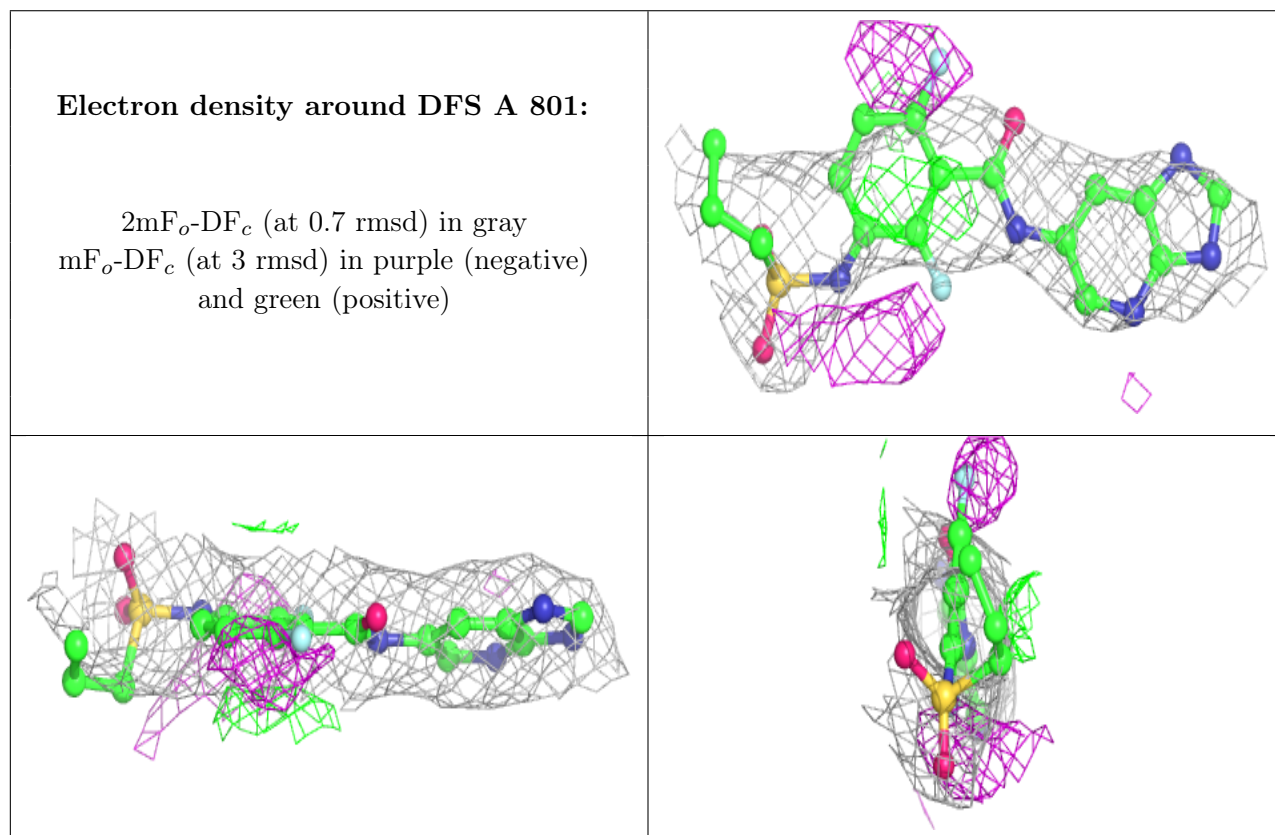
There are no monosaccharides in this entry.

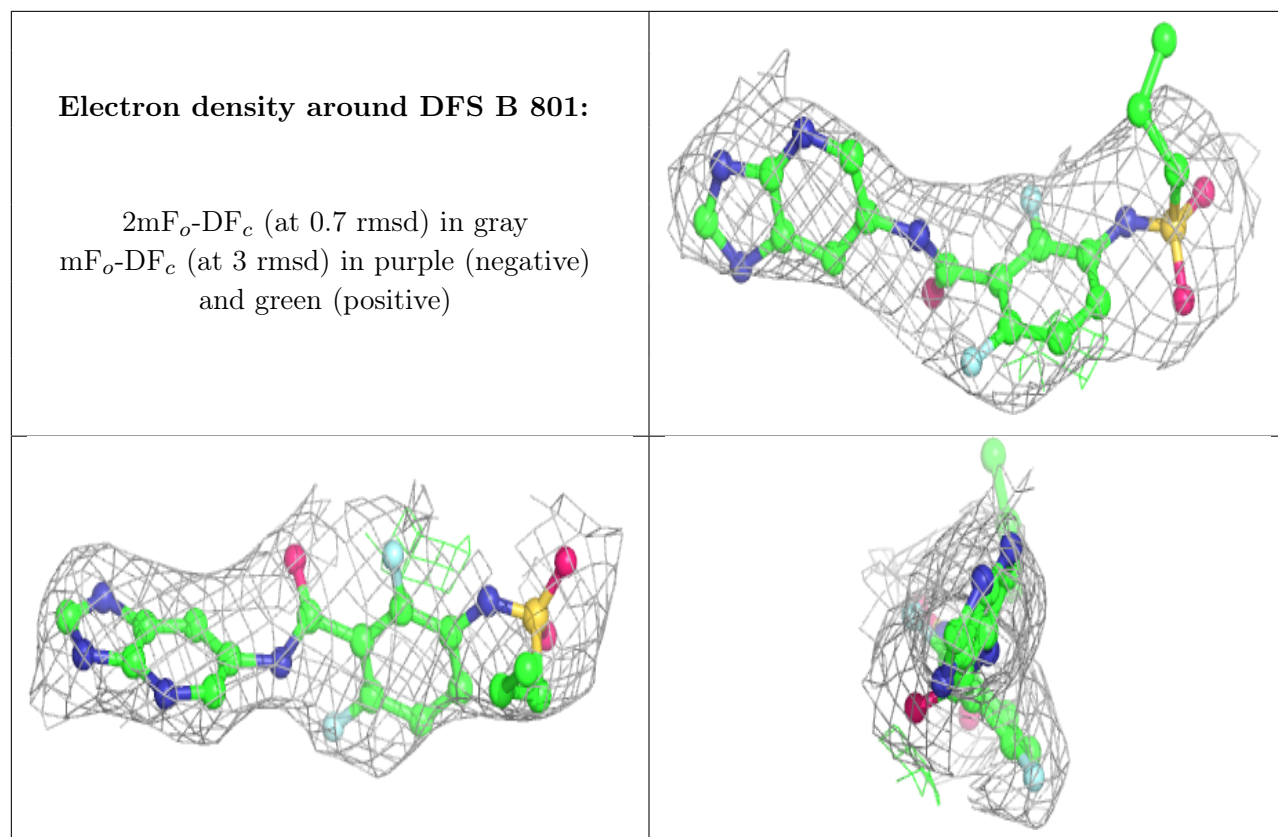
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
2	DFS	A	801	27/27	0.82	0.37	55,71,78,82	0
2	DFS	B	801	27/27	0.93	0.21	29,39,40,43	0

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





6.5 Other polymers [i](#)

There are no such residues in this entry.