



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:20 PM GMT

PDB ID : 4EHG
Title : B-Raf Kinase Domain in Complex with an Aminopyridimine-based Inhibitor
Authors : Voegtli, W.C.
Deposited on : 2012-04-02
Resolution : 3.50 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

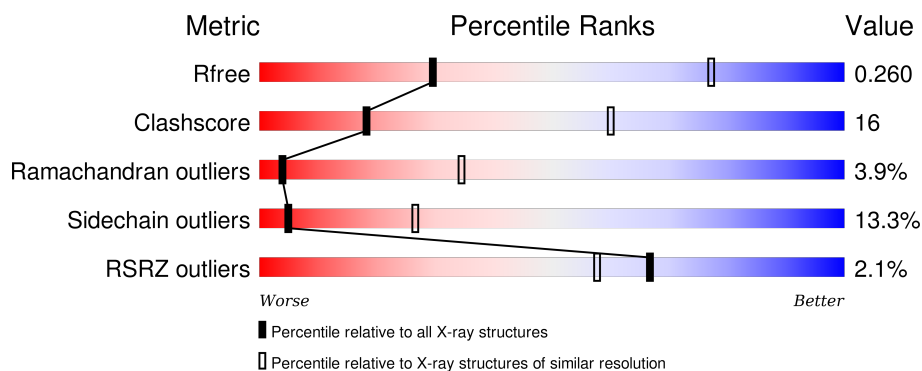
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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}	91344	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	307	<div> <div>54%</div> <div>26%</div> <div>• •</div> <div>15%</div> </div>
1	B	307	<div> <div>55%</div> <div>23%</div> <div>6%</div> <div>•</div> <div>15%</div> </div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 4233 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
1	A	262	Total	C	N	O	S	0	0	0
			2095	1343	364	375	13			
1	B	260	Total	C	N	O	S	0	0	0
			2080	1334	362	371	13			

There are 26 discrepancies between the modelled and reference sequences:

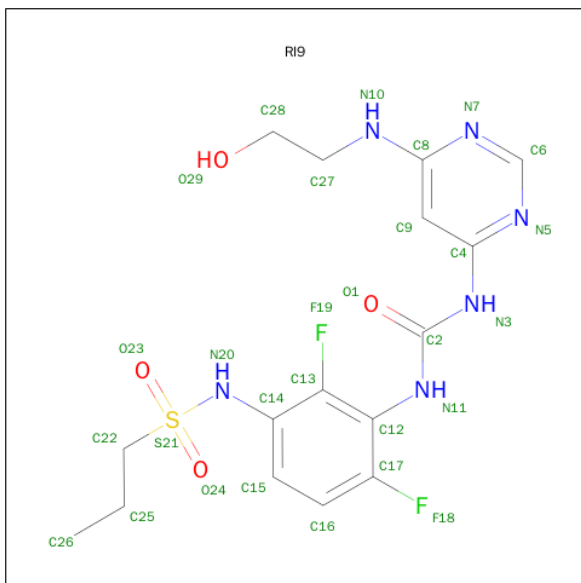
Chain	Residue	Modelled	Actual	Comment	Reference
A	420	MET	-	EXPRESSION TAG	UNP P15056
A	421	ASP	-	EXPRESSION TAG	UNP P15056
A	422	ARG	-	EXPRESSION TAG	UNP P15056
A	423	GLY	-	EXPRESSION TAG	UNP P15056
A	424	SER	-	EXPRESSION TAG	UNP P15056
A	425	HIS	-	EXPRESSION TAG	UNP P15056
A	426	HIS	-	EXPRESSION TAG	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
A	600	GLU	VAL	ENGINEERED MUTATION	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

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Chain	Residue	Modelled	Actual	Comment	Reference
B	600	GLU	VAL	ENGINEERED MUTATION	UNP P15056

- Molecule 2 is N-{2,4-DIFLUORO-3-[(6-[(2-HYDROXYETHYL)AMINO]PYRIMIDIN-4-YL)CARBAMOYL)AMINO]PHENYL}PROPANE-1-SULFONAMIDE (three-letter code: RI9) (formula: C₁₆H₂₀F₂N₆O₄S).

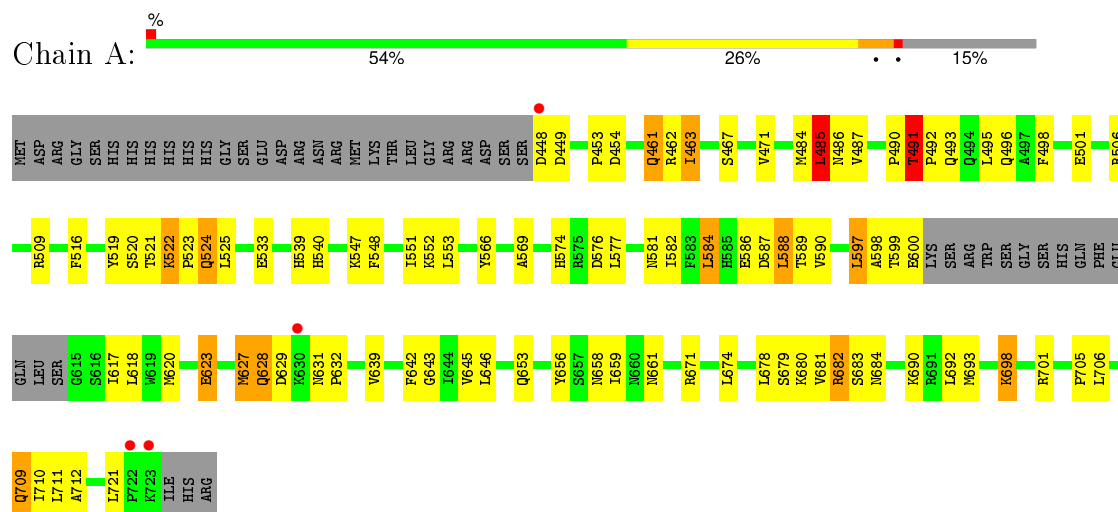


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	0
			29	16	2	6	4	1		
2	B	1	Total	C	F	N	O	S	0	0
			29	16	2	6	4	1		

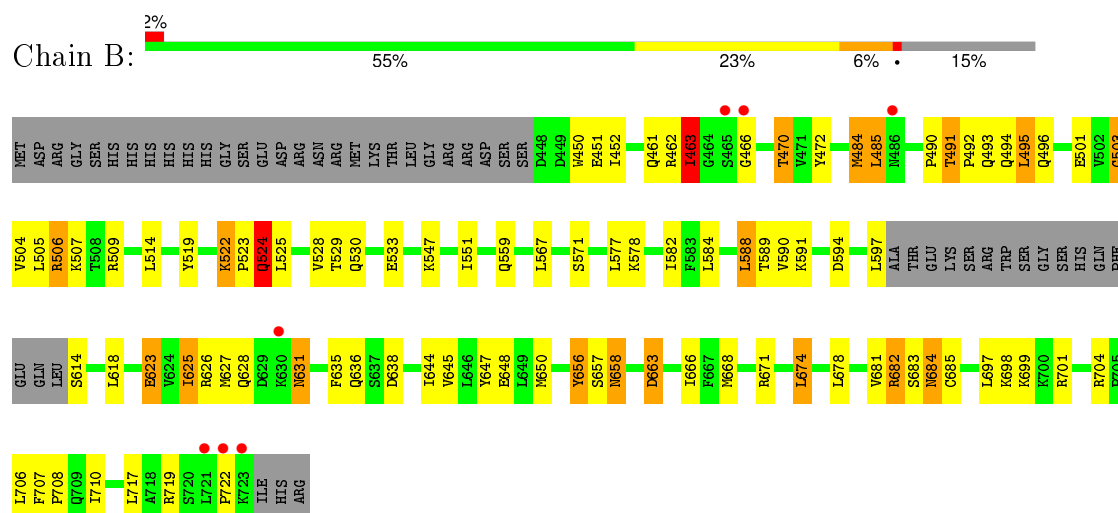
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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	100.50 Å 100.50 Å 161.31 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.57 – 3.50 29.57 – 3.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (29.57-3.50) 97.3 (29.57-3.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.69 (at 3.47 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.212 , 0.270 0.198 , 0.260	Depositor DCC
R_{free} test set	511 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.194	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 31.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 10670 reflections (0.009%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4233	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.57% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: RI9

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	0.57	0/2140	0.73	0/2888
1	B	0.53	0/2125	0.70	2/2867 (0.1%)
All	All	0.55	0/4265	0.72	2/5755 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	495	LEU	CA-CB-CG	6.02	129.15	115.30
1	B	588	LEU	CA-CB-CG	5.21	127.28	115.30

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	2095	0	2132	68	0
1	B	2080	0	2119	63	0
2	A	29	0	20	4	0
2	B	29	0	20	4	0
All	All	4233	0	4291	133	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 16.

All (133) 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:A:801:RI9:C28	2:A:801:RI9:C27	1.78	1.59
1:A:628:GLN:HG3	1:A:629:ASP:H	1.16	1.07
1:A:493:GLN:HB2	1:A:496:GLN:HB3	1.44	0.97
1:B:490:PRO:HA	1:B:491:THR:OG1	1.70	0.91
1:A:599:THR:H	1:A:600:GLU:HA	1.38	0.87
1:A:576:ASP:O	1:A:581:ASN:ND2	2.11	0.84
1:B:668:MET:HB3	1:B:674:LEU:HB2	1.59	0.83
1:A:628:GLN:HG3	1:A:629:ASP:N	1.94	0.79
1:A:582:ILE:HG23	1:A:590:VAL:HG13	1.66	0.78
1:A:521:THR:HA	1:A:525:LEU:HD23	1.64	0.78
1:A:490:PRO:HA	1:A:491:THR:OG1	1.85	0.77
1:B:490:PRO:HA	1:B:491:THR:CB	2.15	0.76
1:B:493:GLN:HB2	1:B:496:GLN:CB	2.17	0.74
1:B:493:GLN:HB2	1:B:496:GLN:HB2	1.69	0.73
1:B:682:ARG:HH11	1:B:684:ASN:HD21	1.36	0.73
1:B:707:PHE:HA	1:B:710:ILE:HB	1.71	0.72
1:A:487:VAL:O	1:A:487:VAL:HG23	1.90	0.71
2:A:801:RI9:N10	2:A:801:RI9:C28	2.54	0.70
1:A:485:LEU:HD11	1:A:597:LEU:HD11	1.74	0.69
1:A:584:LEU:HD13	1:A:587:ASP:HA	1.75	0.68
1:B:658:ASN:H	1:B:658:ASN:HD22	1.42	0.67
1:B:719:ARG:O	1:B:722:PRO:HD3	1.95	0.67
1:B:462:ARG:O	1:B:463:ILE:HG12	1.95	0.67
1:B:684:ASN:H	1:B:684:ASN:HD22	1.43	0.66
1:A:705:PRO:HB2	1:A:709:GLN:HG2	1.78	0.65
1:B:491:THR:N	1:B:492:PRO:HD3	2.12	0.63
1:A:551:ILE:H	1:A:551:ILE:HD12	1.63	0.62
1:A:599:THR:N	1:A:600:GLU:HA	2.13	0.62
1:B:650:MET:HB3	1:B:681:VAL:HG13	1.80	0.61
1:B:623:GLU:OE1	1:B:704:ARG:NH1	2.33	0.61
1:B:505:LEU:HD22	2:B:801:RI9:H22	1.83	0.60
1:A:485:LEU:O	1:A:487:VAL:N	2.35	0.60
1:A:490:PRO:HA	1:A:491:THR:CB	2.34	0.58
1:A:454:ASP:OD1	1:A:523:PRO:HB3	2.03	0.58
1:A:453:PRO:HA	1:A:522:LYS:HD3	1.86	0.58
1:A:681:VAL:HG21	1:A:690:LYS:HD2	1.86	0.58
1:B:503:GLY:HA2	1:B:506:ARG:HD2	1.86	0.58
1:A:493:GLN:HB2	1:A:496:GLN:CB	2.26	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:663:ASP:HA	1:B:666:ILE:HD12	1.87	0.57
2:A:801:RI9:O29	2:A:801:RI9:C27	2.49	0.57
1:A:516:PHE:O	1:B:509:ARG:HD3	2.05	0.56
1:B:551:ILE:HD12	1:B:551:ILE:H	1.70	0.56
1:A:463:ILE:HD11	1:A:471:VAL:HG12	1.86	0.56
1:B:491:THR:N	1:B:492:PRO:CD	2.68	0.55
1:A:709:GLN:O	1:A:712:ALA:HB3	2.07	0.55
1:B:493:GLN:HB2	1:B:496:GLN:HB3	1.89	0.55
1:A:617:ILE:HG22	1:A:620:MET:SD	2.46	0.55
1:B:484:MET:HG3	1:B:524:GLN:HE22	1.72	0.55
1:A:628:GLN:HE21	1:A:629:ASP:CG	2.10	0.55
1:A:461:GLN:NE2	1:A:462:ARG:O	2.40	0.54
1:A:533:GLU:HA	1:A:533:GLU:OE2	2.07	0.54
1:A:639:VAL:O	1:A:642:PHE:HB3	2.08	0.54
1:A:501:GLU:HG3	1:A:597:LEU:HA	1.90	0.53
1:B:491:THR:H	1:B:492:PRO:HD3	1.73	0.52
1:B:514:LEU:HA	1:B:591:LYS:HD2	1.91	0.52
1:B:514:LEU:O	2:B:801:RI9:H26	2.09	0.52
1:A:599:THR:CG2	1:A:600:GLU:C	2.77	0.52
1:A:522:LYS:HB2	1:A:523:PRO:HD3	1.90	0.52
1:A:599:THR:HG22	1:A:600:GLU:C	2.30	0.52
1:B:697:LEU:O	1:B:698:LYS:C	2.48	0.52
1:A:487:VAL:CG2	1:A:487:VAL:O	2.58	0.51
1:B:501:GLU:HG3	1:B:597:LEU:HD23	1.91	0.51
1:A:682:ARG:HD2	1:A:684:ASN:HD21	1.76	0.50
1:A:491:THR:N	1:A:492:PRO:CD	2.74	0.50
1:A:643:GLY:O	1:A:646:LEU:HB2	2.12	0.50
1:B:656:TYR:O	1:B:658:ASN:N	2.45	0.49
1:B:678:LEU:O	1:B:681:VAL:HG23	2.12	0.49
1:B:559:GLN:HB3	1:B:590:VAL:HB	1.94	0.49
1:B:623:GLU:OE1	1:B:701:ARG:HB3	2.12	0.49
1:A:533:GLU:O	1:A:586:GLU:N	2.39	0.49
1:B:451:GLU:HA	1:B:519:TYR:O	2.13	0.48
1:A:574:HIS:C	1:A:576:ASP:H	2.17	0.48
1:A:521:THR:HA	1:A:525:LEU:CD2	2.37	0.48
1:B:707:PHE:O	1:B:708:PRO:C	2.51	0.47
1:B:594:ASP:H	2:B:801:RI9:HN20	1.62	0.47
1:A:678:LEU:O	1:A:680:LYS:N	2.47	0.47
1:A:522:LYS:HB2	1:A:523:PRO:CD	2.44	0.47
1:B:484:MET:HG3	1:B:524:GLN:NE2	2.29	0.47
1:B:636:GLN:N	1:B:636:GLN:OE1	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:548:PHE:HB2	1:A:553:LEU:HD21	1.96	0.46
1:B:635:PHE:O	1:B:638:ASP:HB2	2.16	0.46
1:B:492:PRO:HA	1:B:493:GLN:HA	1.38	0.46
1:A:678:LEU:C	1:A:680:LYS:H	2.19	0.46
1:B:582:ILE:HG23	1:B:590:VAL:HG13	1.97	0.46
1:A:627:MET:HA	1:A:627:MET:CE	2.46	0.46
1:A:706:LEU:O	1:A:710:ILE:HG13	2.16	0.46
1:B:452:ILE:HD12	1:B:528:VAL:HG21	1.98	0.46
1:A:551:ILE:N	1:A:551:ILE:HD12	2.30	0.46
2:A:801:RI9:HO29	2:A:801:RI9:C27	2.27	0.45
1:B:682:ARG:O	1:B:684:ASN:N	2.49	0.45
1:B:504:VAL:O	1:B:507:LYS:HB2	2.16	0.45
1:B:522:LYS:O	1:B:523:PRO:C	2.54	0.45
1:B:485:LEU:HD23	1:B:525:LEU:HB3	1.99	0.45
1:A:698:LYS:HA	1:A:698:LYS:HD2	1.77	0.45
1:B:493:GLN:CB	1:B:496:GLN:HB2	2.42	0.45
1:A:448:ASP:OD2	1:A:448:ASP:C	2.54	0.45
1:A:498:PHE:CD1	1:A:525:LEU:HD13	2.52	0.45
1:B:529:THR:HG21	2:B:801:RI9:H25	1.98	0.45
1:B:656:TYR:N	1:B:656:TYR:CD1	2.84	0.44
1:A:692:LEU:HD11	1:A:710:ILE:HG23	2.00	0.44
1:B:645:VAL:HA	1:B:648:GLU:HG3	2.00	0.44
1:B:681:VAL:HG12	1:B:685:CYS:HB3	1.98	0.44
1:A:584:LEU:HD23	1:A:584:LEU:HA	1.75	0.44
1:A:628:GLN:CG	1:A:629:ASP:H	2.05	0.44
1:B:644:ILE:O	1:B:647:TYR:HB3	2.18	0.44
1:A:519:TYR:C	1:A:519:TYR:CD1	2.91	0.44
1:B:491:THR:H	1:B:492:PRO:CD	2.28	0.43
1:B:462:ARG:C	1:B:463:ILE:HG12	2.37	0.43
1:B:530:GLN:HG2	1:B:530:GLN:O	2.18	0.43
1:B:551:ILE:N	1:B:551:ILE:HD12	2.33	0.43
1:A:646:LEU:HB3	1:A:693:MET:HG3	2.00	0.43
1:A:539:HIS:O	1:A:540:HIS:C	2.55	0.43
1:B:656:TYR:C	1:B:658:ASN:H	2.22	0.43
1:A:681:VAL:O	1:A:681:VAL:HG23	2.18	0.43
1:A:678:LEU:C	1:A:680:LYS:N	2.72	0.43
1:A:659:ILE:HG22	1:A:661:ASN:H	1.84	0.43
1:B:470:THR:HG22	1:B:472:TYR:CE1	2.54	0.42
1:A:711:LEU:O	1:A:712:ALA:C	2.57	0.42
1:B:644:ILE:O	1:B:648:GLU:HG3	2.19	0.42
1:A:509:ARG:HD2	1:B:450:TRP:CH2	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:490:PRO:CA	1:B:491:THR:CB	2.95	0.42
1:B:491:THR:HG21	1:B:494:GLN:OE1	2.20	0.42
1:B:567:LEU:HD23	1:B:567:LEU:HA	1.92	0.42
1:A:597:LEU:O	1:A:599:THR:N	2.52	0.42
1:A:520:SER:O	1:A:525:LEU:HA	2.20	0.41
1:A:492:PRO:HA	1:A:493:GLN:HA	1.33	0.41
1:A:523:PRO:HG2	1:A:524:GLN:H	1.85	0.41
1:A:566:TYR:O	1:A:569:ALA:HB3	2.20	0.41
1:B:631:ASN:HD22	1:B:631:ASN:HA	1.61	0.41
1:A:599:THR:HG23	1:A:600:GLU:C	2.42	0.40
1:A:623:GLU:HG3	1:A:623:GLU:H	1.65	0.40
1:B:625:ILE:HG22	1:B:626:ARG:N	2.35	0.40
1:A:631:ASN:HA	1:A:632:PRO:HD3	1.66	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	258/307 (84%)	216 (84%)	31 (12%)	11 (4%)	3	31
1	B	256/307 (83%)	220 (86%)	27 (10%)	9 (4%)	4	38
All	All	514/614 (84%)	436 (85%)	58 (11%)	20 (4%)	4	34

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	486	ASN
1	A	598	ALA
1	B	491	THR
1	B	683	SER
1	A	628	GLN

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Mol	Chain	Res	Type
1	A	683	SER
1	B	461	GLN
1	B	466	GLY
1	B	524	GLN
1	B	657	SER
1	A	467	SER
1	A	491	THR
1	A	588	LEU
1	A	597	LEU
1	A	679	SER
1	A	721	LEU
1	A	485	LEU
1	B	463	ILE
1	B	503	GLY
1	B	625	ILE

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	230/271 (85%)	201 (87%)	29 (13%)	5	28
1	B	229/271 (84%)	197 (86%)	32 (14%)	4	24
All	All	459/542 (85%)	398 (87%)	61 (13%)	5	26

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	449	ASP
1	A	461	GLN
1	A	463	ILE
1	A	484	MET
1	A	485	LEU
1	A	491	THR
1	A	495	LEU
1	A	506	ARG

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Mol	Chain	Res	Type
1	A	522	LYS
1	A	524	GLN
1	A	547	LYS
1	A	552	LYS
1	A	577	LEU
1	A	584	LEU
1	A	588	LEU
1	A	589	THR
1	A	618	LEU
1	A	623	GLU
1	A	627	MET
1	A	645	VAL
1	A	653	GLN
1	A	656	TYR
1	A	658	ASN
1	A	671	ARG
1	A	674	LEU
1	A	682	ARG
1	A	698	LYS
1	A	701	ARG
1	A	709	GLN
1	B	463	ILE
1	B	470	THR
1	B	484	MET
1	B	485	LEU
1	B	495	LEU
1	B	506	ARG
1	B	522	LYS
1	B	524	GLN
1	B	533	GLU
1	B	547	LYS
1	B	571	SER
1	B	577	LEU
1	B	578	LYS
1	B	584	LEU
1	B	588	LEU
1	B	589	THR
1	B	614	SER
1	B	618	LEU
1	B	623	GLU
1	B	627	MET
1	B	628	GLN

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Mol	Chain	Res	Type
1	B	631	ASN
1	B	656	TYR
1	B	658	ASN
1	B	663	ASP
1	B	671	ARG
1	B	674	LEU
1	B	682	ARG
1	B	684	ASN
1	B	699	LYS
1	B	706	LEU
1	B	717	LEU

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

Mol	Chain	Res	Type
1	A	461	GLN
1	A	486	ASN
1	A	562	GLN
1	A	628	GLN
1	A	658	ASN
1	A	684	ASN
1	B	456	GLN
1	B	486	ASN
1	B	539	HIS
1	B	562	GLN
1	B	585	HIS
1	B	631	ASN
1	B	653	GLN
1	B	658	ASN
1	B	684	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	RI9	A	801	-	30,30,30	2.41	10 (33%)	38,41,41	1.93	10 (26%)
2	RI9	B	801	-	30,30,30	1.72	6 (20%)	38,41,41	1.98	12 (31%)

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	RI9	A	801	-	-	0/21/21/21	0/2/2/2
2	RI9	B	801	-	-	0/21/21/21	0/2/2/2

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	RI9	C14-N20	-5.20	1.33	1.42
2	A	801	RI9	C14-N20	-4.85	1.34	1.42
2	A	801	RI9	C2-N11	-2.14	1.33	1.37
2	A	801	RI9	C2-N3	-2.06	1.33	1.37
2	A	801	RI9	O24-S21	2.13	1.46	1.43
2	B	801	RI9	C22-S21	2.18	1.84	1.78
2	A	801	RI9	O23-S21	2.28	1.47	1.43
2	A	801	RI9	C22-S21	2.31	1.84	1.78
2	B	801	RI9	O24-S21	2.37	1.47	1.43
2	A	801	RI9	C12-C17	2.58	1.41	1.38
2	B	801	RI9	O23-S21	2.65	1.47	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	RI9	C12-C17	2.84	1.41	1.38
2	B	801	RI9	S21-N20	3.55	1.71	1.62
2	A	801	RI9	S21-N20	4.02	1.73	1.62
2	A	801	RI9	C27-N10	5.15	1.58	1.46
2	A	801	RI9	C27-C28	7.68	1.78	1.50

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	RI9	C28-C27-N10	-5.08	98.13	111.65
2	B	801	RI9	C12-C13-C14	-3.96	119.50	124.47
2	A	801	RI9	O29-C28-C27	-3.52	97.74	111.64
2	B	801	RI9	C13-C12-N11	-3.38	115.57	121.46
2	A	801	RI9	C9-C8-N7	-3.29	119.38	123.06
2	B	801	RI9	C9-C4-N5	-3.25	119.43	123.06
2	A	801	RI9	C12-C13-C14	-3.10	120.58	124.47
2	B	801	RI9	C16-C17-C12	-2.76	119.41	122.81
2	A	801	RI9	C9-C4-N3	-2.67	115.42	122.81
2	A	801	RI9	C9-C4-N5	-2.64	120.10	123.06
2	B	801	RI9	C22-S21-N20	-2.57	97.13	105.13
2	B	801	RI9	C9-C4-N3	-2.39	116.17	122.81
2	B	801	RI9	C4-N3-C2	-2.16	127.53	130.53
2	B	801	RI9	O23-S21-C22	2.38	112.22	107.73
2	B	801	RI9	C12-N11-C2	2.64	125.93	121.90
2	A	801	RI9	F18-C17-C12	2.93	120.26	117.74
2	B	801	RI9	F18-C17-C12	3.13	120.44	117.74
2	A	801	RI9	C17-C12-C13	3.24	119.14	115.60
2	B	801	RI9	N3-C4-N5	3.32	124.40	114.83
2	A	801	RI9	N3-C4-N5	3.35	124.47	114.83
2	A	801	RI9	C12-N11-C2	3.95	127.94	121.90
2	B	801	RI9	C17-C12-C13	4.98	121.03	115.60

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	RI9	4	0
2	B	801	RI9	4	0

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	262/307 (85%)	-0.50	4 (1%) 76 67	10, 24, 51, 77	0
1	B	260/307 (84%)	-0.32	7 (2%) 58 47	7, 27, 62, 70	0
All	All	522/614 (85%)	-0.41	11 (2%) 67 58	7, 26, 57, 77	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	723	LYS	3.7
1	A	723	LYS	3.1
1	A	448	ASP	3.0
1	B	466	GLY	2.8
1	B	722	PRO	2.6
1	B	721	LEU	2.4
1	B	630	LYS	2.4
1	B	465	SER	2.2
1	A	722	PRO	2.1
1	A	630	LYS	2.1
1	B	486	ASN	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	RI9	A	801	29/29	0.95	0.20	0.74	33,36,38,40	0
2	RI9	B	801	29/29	0.96	0.20	0.12	41,44,47,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.