



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:02 AM GMT

PDB ID : 3D4Q  
Title : Pyrazole-based inhibitors of B-Raf kinase  
Authors : Morales, T.; Vigers, G.P.A; Brandhuber, B.J  
Deposited on : 2008-05-14  
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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

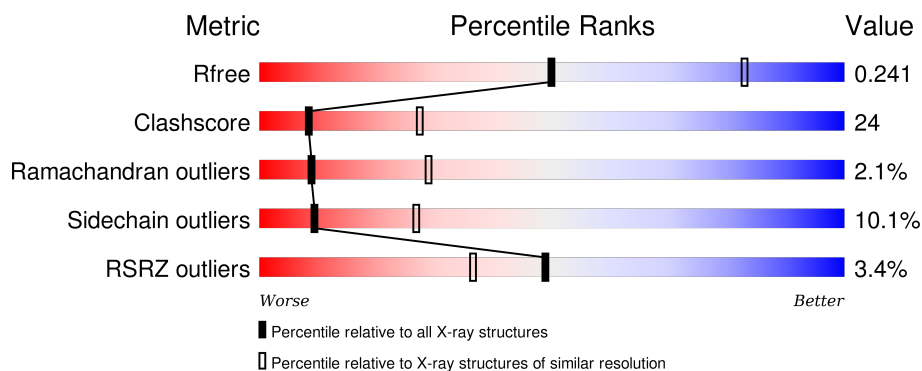
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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>2%</div> <div>56%</div> <div>24%</div> <div>5%</div> <div>14%</div> </div>
1	B	307	<div> <div>4%</div> <div>55%</div> <div>24%</div> <div>6%</div> <div>14%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4293 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 B-Raf proto-oncogene serine/threonine-protein kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2107	1352	366	376	13			
1	B	264	Total	C	N	O	S	0	0	0
			2107	1352	366	376	13			

There are 26 discrepancies between the modelled and reference sequences:

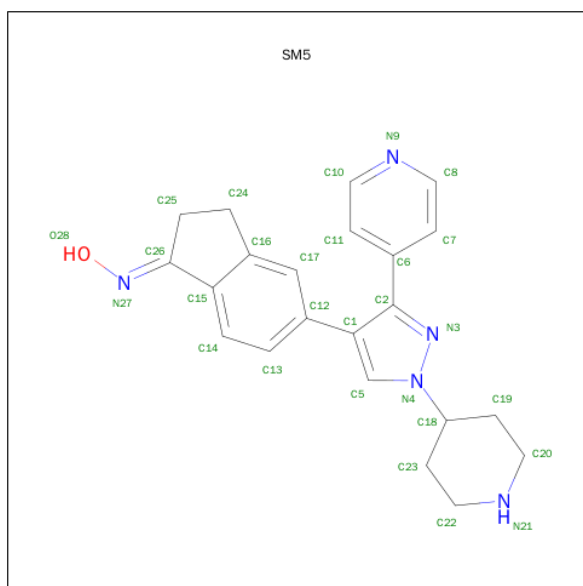
Chain	Residue	Modelled	Actual	Comment	Reference
A	420	MET	-	EXPRESSION TAG	UNP P15056
A	421	ASP	-	EXPRESSION TAG	UNP P15056
A	422	ARG	-	EXPRESSION TAG	UNP P15056
A	423	GLY	-	EXPRESSION TAG	UNP P15056
A	424	SER	-	EXPRESSION TAG	UNP P15056
A	425	HIS	-	EXPRESSION TAG	UNP P15056
A	426	HIS	-	EXPRESSION TAG	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	432	SER	-	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

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Chain	Residue	Modelled	Actual	Comment	Reference
B	432	SER	-	EXPRESSION TAG	UNP P15056

- Molecule 2 is (1E)-5-(1-PIPERIDIN-4-YL-3-PYRIDIN-4-YL-1H-PYRAZOL-4-YL)-2,3-DIHYDRO-1H-INDEN-1-ONE OXIME (three-letter code: SM5) (formula: C<sub>22</sub>H<sub>23</sub>N<sub>5</sub>O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			28	22	5	1		
2	B	1	Total	C	N	O	0	0
			28	22	5	1		

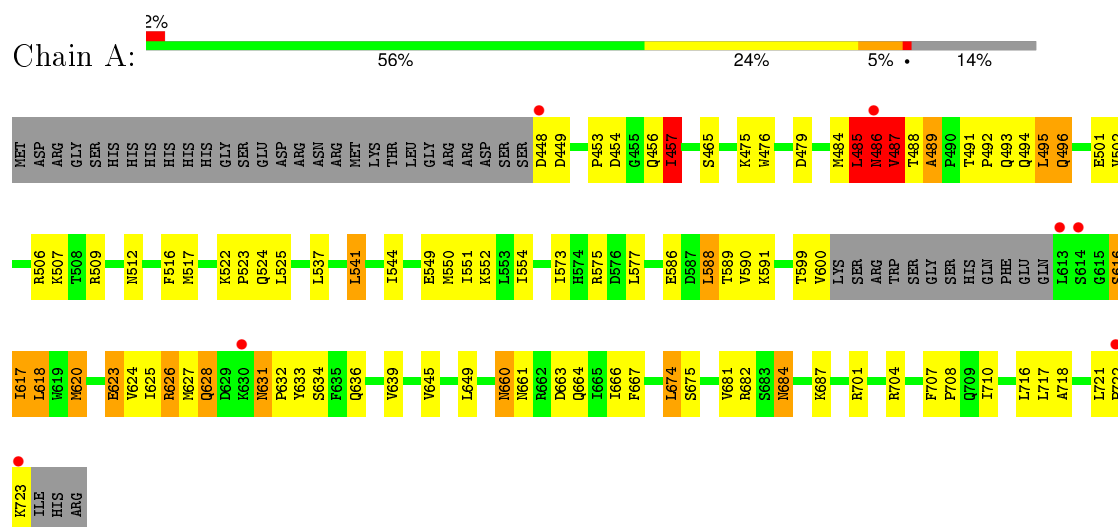
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	B	14	Total	O	0	0
			14	14		

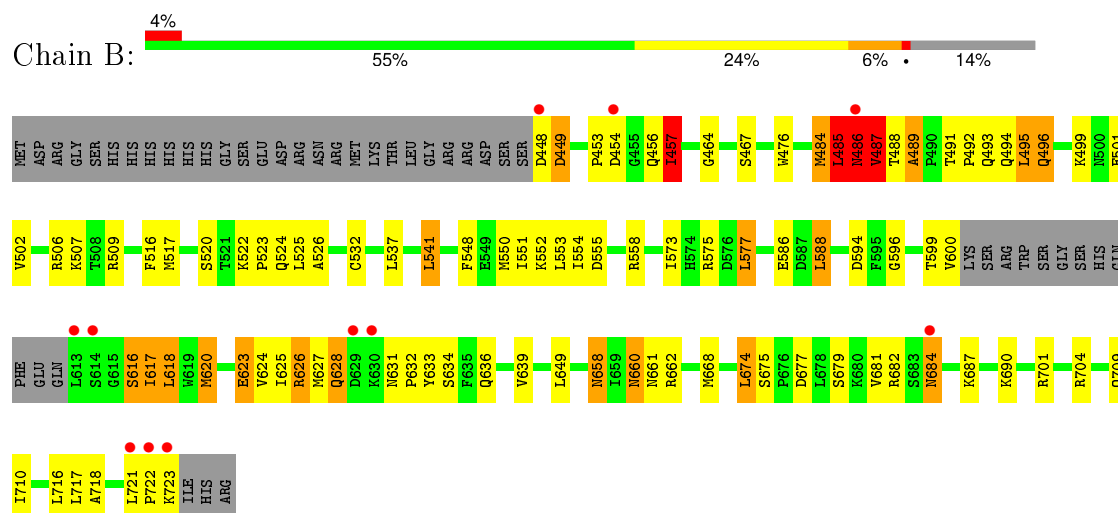
### 3 Residue-property plots

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: B-Raf proto-oncogene serine/threonine-protein kinase



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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	100.34Å 100.34Å 162.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 47.94 – 2.81	Depositor EDS
% Data completeness (in resolution range)	97.8 (30.00-2.80) 99.1 (47.94-2.81)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.46 (at 2.81Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.216 , 0.250 0.214 , 0.241	Depositor DCC
$R_{free}$ test set	1002 reflections (4.85%)	DCC
Wilson B-factor (Å <sup>2</sup> )	65.4	Xtriage
Anisotropy	0.229	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 33.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 20661 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4293	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SM5

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	0.70	1/2152 (0.0%)	0.77	2/2905 (0.1%)
1	B	0.69	0/2152	0.76	2/2905 (0.1%)
All	All	0.70	1/4304 (0.0%)	0.76	4/5810 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	645	VAL	CA-CB	5.74	1.66	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	485	LEU	CA-CB-CG	5.78	128.59	115.30
1	A	457	ILE	CG1-CB-CG2	-5.58	99.12	111.40
1	B	485	LEU	CA-CB-CG	5.50	127.94	115.30
1	B	457	ILE	CG1-CB-CG2	-5.04	100.31	111.40

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	485	LEU	Peptide
1	A	486	ASN	Peptide
1	B	485	LEU	Peptide
1	B	486	ASN	Peptide

## 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2107	0	2151	103	0
1	B	2107	0	2151	103	0
2	A	28	0	23	6	0
2	B	28	0	23	8	0
3	A	9	0	0	0	0
3	B	14	0	0	0	0
All	All	4293	0	4348	208	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 24.

All (208) 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:517:MET:SD	1:A:517:MET:CE	2.01	1.48
1:B:517:MET:CE	1:B:517:MET:SD	2.03	1.47
2:B:1:SM5:N3	2:B:1:SM5:C2	1.75	1.31
2:A:1:SM5:C2	2:A:1:SM5:N3	1.68	1.30
1:A:588:LEU:HD21	1:B:586:GLU:HG3	1.39	1.02
1:B:491:THR:HG22	1:B:493:GLN:H	1.23	1.00
1:A:586:GLU:HG3	1:B:588:LEU:HD21	1.41	0.99
1:A:491:THR:HG22	1:A:493:GLN:H	1.24	0.98
1:A:551:ILE:HD12	1:A:551:ILE:H	1.28	0.95
1:B:551:ILE:H	1:B:551:ILE:HD12	1.31	0.95
1:B:668:MET:HB3	1:B:674:LEU:HB2	1.63	0.81
1:A:684:ASN:HD22	1:A:684:ASN:H	1.28	0.81
1:A:554:ILE:HD12	1:A:721:LEU:HD22	1.63	0.80
1:B:554:ILE:HD12	1:B:721:LEU:HD22	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:491:THR:HB	1:A:494:GLN:HG3	1.68	0.76
1:B:454:ASP:HB2	1:B:522:LYS:NZ	2.00	0.76
1:B:491:THR:HB	1:B:494:GLN:HG3	1.66	0.75
1:B:684:ASN:H	1:B:684:ASN:HD22	1.30	0.75
1:A:663:ASP:HA	1:A:666:ILE:HG12	1.68	0.74
1:B:658:ASN:HD22	1:B:658:ASN:H	1.37	0.73
1:A:537:LEU:HD21	1:A:649:LEU:HD21	1.69	0.73
1:B:660:ASN:HD22	1:B:660:ASN:C	1.92	0.72
1:A:617:ILE:HA	1:A:620:MET:HB2	1.71	0.72
1:B:491:THR:CG2	1:B:492:PRO:HD2	2.20	0.72
1:B:617:ILE:HG13	1:B:618:LEU:N	2.05	0.71
1:A:666:ILE:HG13	1:A:667:PHE:N	2.06	0.71
1:A:501:GLU:OE1	2:A:1:SM5:O28	2.08	0.71
1:B:617:ILE:HA	1:B:620:MET:HB2	1.72	0.70
1:B:551:ILE:CD1	1:B:551:ILE:H	2.00	0.70
1:B:537:LEU:HD21	1:B:649:LEU:HD21	1.75	0.69
1:B:616:SER:O	1:B:617:ILE:HG12	1.93	0.68
1:A:486:ASN:O	1:A:487:VAL:HG13	1.92	0.68
1:B:454:ASP:HB2	1:B:522:LYS:HZ3	1.57	0.68
1:B:486:ASN:O	1:B:487:VAL:HG13	1.94	0.67
1:A:491:THR:CG2	1:A:492:PRO:HD2	2.24	0.67
1:A:551:ILE:CD1	1:A:551:ILE:H	2.00	0.66
1:B:486:ASN:HD22	1:B:487:VAL:HA	1.60	0.66
1:B:487:VAL:O	1:B:489:ALA:N	2.25	0.66
1:B:454:ASP:H	1:B:522:LYS:NZ	1.95	0.65
1:B:684:ASN:N	1:B:684:ASN:HD22	1.90	0.64
1:A:626:ARG:O	1:A:628:GLN:N	2.30	0.64
1:A:551:ILE:N	1:A:551:ILE:HD12	2.09	0.64
1:A:718:ALA:O	1:A:722:PRO:HD3	1.99	0.63
1:A:684:ASN:N	1:A:684:ASN:HD22	1.90	0.62
1:B:718:ALA:O	1:B:722:PRO:HD3	1.99	0.62
1:B:464:GLY:HA3	2:B:1:SM5:H119	1.82	0.62
2:A:1:SM5:H11	2:A:1:SM5:C17	2.30	0.62
1:A:684:ASN:ND2	1:A:684:ASN:H	1.96	0.62
1:A:486:ASN:HD22	1:A:487:VAL:HA	1.65	0.62
1:A:661:ASN:OD1	1:A:664:GLN:HG2	1.99	0.62
1:A:454:ASP:HB2	1:A:522:LYS:HZ3	1.64	0.61
1:B:573:ILE:HG22	1:B:575:ARG:HG3	1.82	0.61
1:A:660:ASN:C	1:A:660:ASN:HD22	2.03	0.61
1:B:491:THR:HG23	1:B:492:PRO:HD2	1.82	0.61
1:A:666:ILE:HG13	1:A:667:PHE:H	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:684:ASN:H	1:B:684:ASN:ND2	1.99	0.60
1:B:660:ASN:ND2	1:B:660:ASN:C	2.53	0.59
1:A:616:SER:O	1:A:617:ILE:HG12	2.02	0.59
1:A:682:ARG:HG3	1:A:684:ASN:ND2	2.16	0.59
1:A:523:PRO:O	1:A:524:GLN:HG3	2.03	0.59
1:B:682:ARG:HG3	1:B:684:ASN:ND2	2.17	0.59
2:B:1:SM5:C17	2:B:1:SM5:H11	2.33	0.59
1:B:626:ARG:O	1:B:628:GLN:N	2.35	0.59
1:A:617:ILE:HG13	1:A:618:LEU:N	2.18	0.58
1:A:491:THR:HG23	1:A:492:PRO:HD2	1.85	0.58
1:A:454:ASP:H	1:A:522:LYS:NZ	2.01	0.58
1:A:599:THR:HG22	1:A:600:VAL:N	2.19	0.57
2:B:1:SM5:H11	2:B:1:SM5:C12	2.35	0.56
1:B:586:GLU:O	1:B:588:LEU:HD12	2.05	0.56
1:B:491:THR:HG22	1:B:492:PRO:CD	2.36	0.55
1:B:620:MET:HE2	1:B:625:ILE:HA	1.88	0.55
1:B:486:ASN:ND2	1:B:486:ASN:C	2.60	0.55
1:A:488:THR:O	1:A:489:ALA:HB2	2.07	0.55
2:A:1:SM5:H11	2:A:1:SM5:C12	2.36	0.55
1:A:663:ASP:O	1:A:666:ILE:CG1	2.55	0.55
1:A:465:SER:H	2:A:1:SM5:H120	1.71	0.54
1:B:491:THR:CG2	1:B:493:GLN:H	2.09	0.54
1:B:620:MET:CE	1:B:625:ILE:HA	2.37	0.54
1:B:454:ASP:CB	1:B:522:LYS:NZ	2.68	0.54
1:B:488:THR:O	1:B:489:ALA:HB2	2.07	0.54
1:B:717:LEU:HD12	1:B:717:LEU:O	2.07	0.54
1:A:681:VAL:O	1:A:681:VAL:HG23	2.07	0.54
1:A:620:MET:HG2	1:A:624:VAL:CG1	2.38	0.54
1:B:716:LEU:HD23	1:B:716:LEU:C	2.29	0.53
1:A:717:LEU:C	1:A:717:LEU:HD12	2.29	0.53
1:B:486:ASN:CG	1:B:524:GLN:HB3	2.29	0.53
1:A:488:THR:O	1:A:489:ALA:CB	2.57	0.52
1:A:487:VAL:O	1:A:489:ALA:N	2.37	0.52
1:B:491:THR:HG22	1:B:492:PRO:HD2	1.88	0.52
1:A:586:GLU:O	1:A:588:LEU:HD12	2.09	0.52
1:A:486:ASN:C	1:A:486:ASN:ND2	2.62	0.52
1:A:624:VAL:HA	1:A:632:PRO:HB2	1.92	0.52
1:B:502:VAL:O	1:B:506:ARG:HG2	2.09	0.52
1:A:663:ASP:O	1:A:666:ILE:HG13	2.10	0.52
1:A:486:ASN:CG	1:A:524:GLN:HB3	2.30	0.52
1:B:501:GLU:OE1	2:B:1:SM5:O28	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:550:MET:O	1:B:554:ILE:HG13	2.10	0.51
1:B:658:ASN:HD22	1:B:658:ASN:N	2.05	0.51
1:B:624:VAL:HA	1:B:632:PRO:HB2	1.92	0.51
1:B:620:MET:HG2	1:B:624:VAL:HG12	1.91	0.51
1:A:484:MET:HB3	1:A:486:ASN:OD1	2.10	0.51
1:B:620:MET:HG2	1:B:624:VAL:CG1	2.41	0.51
1:B:626:ARG:O	1:B:626:ARG:HG3	2.11	0.51
1:A:674:LEU:HD22	1:A:675:SER:N	2.26	0.50
1:B:717:LEU:C	1:B:717:LEU:HD12	2.31	0.50
1:B:599:THR:HG22	1:B:600:VAL:N	2.26	0.50
1:A:496:GLN:OE1	1:A:496:GLN:HA	2.10	0.50
1:B:660:ASN:HD22	1:B:661:ASN:N	2.10	0.50
1:B:532:CYS:H	2:B:1:SM5:C8	2.24	0.50
1:B:537:LEU:HG	1:B:541:LEU:HD22	1.94	0.50
1:A:454:ASP:HB2	1:A:522:LYS:NZ	2.26	0.50
1:B:457:ILE:HG22	1:B:476:TRP:HB2	1.94	0.49
1:B:454:ASP:CB	1:B:522:LYS:HZ1	2.25	0.49
1:B:488:THR:O	1:B:489:ALA:CB	2.60	0.49
1:B:495:LEU:HD22	1:B:525:LEU:HD11	1.95	0.49
1:A:491:THR:HG22	1:A:493:GLN:N	2.09	0.49
1:A:491:THR:HG22	1:A:492:PRO:CD	2.43	0.49
1:A:507:LYS:HG2	1:B:448:ASP:HA	1.94	0.49
1:B:491:THR:HG21	1:B:493:GLN:HB3	1.95	0.48
1:A:626:ARG:O	1:A:626:ARG:HG3	2.13	0.48
1:A:495:LEU:HD22	1:A:525:LEU:HD11	1.94	0.48
1:A:716:LEU:C	1:A:716:LEU:HD23	2.33	0.48
1:A:554:ILE:HD12	1:A:721:LEU:CD2	2.39	0.48
1:B:684:ASN:N	1:B:684:ASN:ND2	2.60	0.48
1:A:620:MET:HE2	1:A:625:ILE:HA	1.94	0.48
1:B:575:ARG:HG2	1:B:633:TYR:CD2	2.49	0.48
1:A:453:PRO:HG2	1:A:456:GLN:NE2	2.28	0.48
1:A:575:ARG:HG2	1:A:633:TYR:CD2	2.47	0.48
1:B:681:VAL:O	1:B:681:VAL:HG23	2.13	0.48
1:B:484:MET:HB3	1:B:486:ASN:OD1	2.15	0.47
1:B:551:ILE:N	1:B:551:ILE:HD12	2.14	0.47
1:A:684:ASN:N	1:A:684:ASN:ND2	2.59	0.47
2:A:1:SM5:C11	2:A:1:SM5:C17	2.93	0.47
1:A:454:ASP:H	1:A:522:LYS:HZ3	1.63	0.47
1:B:467:SER:OG	1:B:596:GLY:HA3	2.15	0.47
1:A:509:ARG:HD3	1:B:516:PHE:O	2.15	0.47
1:B:520:SER:HB3	1:B:526:ALA:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:MET:HG2	1:A:624:VAL:HG12	1.96	0.47
1:A:457:ILE:HG22	1:A:476:TRP:HB2	1.97	0.47
1:A:623:GLU:HG3	1:A:623:GLU:H	1.22	0.46
1:B:491:THR:CG2	1:B:492:PRO:CD	2.91	0.46
1:B:491:THR:HG22	1:B:492:PRO:N	2.31	0.46
1:A:502:VAL:O	1:A:506:ARG:HG2	2.15	0.46
1:A:491:THR:HG21	1:A:493:GLN:HB3	1.97	0.46
1:A:589:THR:HG22	1:A:591:LYS:HE2	1.98	0.46
1:B:491:THR:HG22	1:B:493:GLN:N	2.08	0.45
1:B:623:GLU:OE1	1:B:704:ARG:NH1	2.47	0.45
1:A:623:GLU:OE1	1:A:704:ARG:NH1	2.46	0.45
1:A:663:ASP:O	1:A:666:ILE:HG12	2.16	0.45
1:A:551:ILE:CD1	1:A:551:ILE:N	2.77	0.45
1:A:631:ASN:HA	1:A:632:PRO:HD2	1.80	0.44
1:B:588:LEU:CD1	1:B:588:LEU:H	2.30	0.44
1:B:541:LEU:HD12	1:B:541:LEU:HA	1.86	0.44
1:A:523:PRO:HB2	1:A:524:GLN:HE21	1.82	0.44
1:B:453:PRO:HG2	1:B:456:GLN:NE2	2.32	0.44
1:A:491:THR:CG2	1:A:493:GLN:H	2.12	0.44
1:B:523:PRO:O	1:B:524:GLN:HG3	2.18	0.43
1:A:636:GLN:HG2	1:A:701:ARG:O	2.18	0.43
1:B:491:THR:CG2	1:B:493:GLN:HB3	2.48	0.43
1:B:636:GLN:HG2	1:B:701:ARG:O	2.19	0.43
1:B:639:VAL:HG13	1:B:710:ILE:HD11	2.00	0.43
1:A:475:LYS:HE3	1:A:479:ASP:OD1	2.18	0.43
1:B:555:ASP:OD1	1:B:558:ARG:NH1	2.51	0.43
1:A:620:MET:CE	1:A:625:ILE:HA	2.48	0.43
1:B:448:ASP:O	1:B:449:ASP:O	2.36	0.43
1:B:499:LYS:HD3	1:B:499:LYS:HA	1.72	0.43
1:A:549:GLU:HB3	1:A:551:ILE:HD13	2.01	0.43
1:A:717:LEU:O	1:A:717:LEU:HD12	2.18	0.43
1:B:554:ILE:HD13	1:B:717:LEU:HD11	2.00	0.43
1:B:548:PHE:HB2	1:B:553:LEU:HD21	2.00	0.43
1:B:486:ASN:HD22	1:B:487:VAL:CA	2.27	0.43
1:A:454:ASP:CB	1:A:522:LYS:NZ	2.82	0.43
1:B:677:ASP:OD1	1:B:679:SER:HB2	2.18	0.43
1:A:491:THR:CG2	1:A:492:PRO:CD	2.95	0.42
1:A:523:PRO:C	1:A:524:GLN:HG3	2.40	0.42
1:B:634:SER:OG	1:B:636:GLN:HB2	2.19	0.42
1:A:516:PHE:O	1:B:509:ARG:HD3	2.19	0.42
1:A:491:THR:CG2	1:A:493:GLN:HB3	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:634:SER:OG	1:A:636:GLN:HB2	2.20	0.42
1:A:512:ASN:HA	1:A:590:VAL:O	2.20	0.42
1:A:721:LEU:C	1:A:723:LYS:N	2.73	0.42
1:B:721:LEU:C	1:B:723:LYS:N	2.73	0.42
1:B:616:SER:O	1:B:617:ILE:CG1	2.65	0.42
1:A:453:PRO:O	1:A:456:GLN:HG3	2.20	0.42
1:A:588:LEU:H	1:A:588:LEU:CD1	2.33	0.42
1:A:707:PHE:N	1:A:708:PRO:CD	2.83	0.42
1:A:448:ASP:HA	1:B:507:LYS:HG2	2.01	0.41
1:B:457:ILE:HG23	1:B:457:ILE:HD13	1.82	0.41
1:B:681:VAL:HG22	1:B:690:LYS:NZ	2.36	0.41
1:A:660:ASN:C	1:A:660:ASN:ND2	2.72	0.41
1:A:682:ARG:HG3	1:A:684:ASN:HD21	1.84	0.41
1:B:496:GLN:OE1	1:B:496:GLN:HA	2.19	0.41
2:B:1:SM5:C17	2:B:1:SM5:C11	2.99	0.41
1:A:674:LEU:HD22	1:A:675:SER:H	1.85	0.41
1:B:464:GLY:HA3	2:B:1:SM5:C19	2.49	0.41
1:A:537:LEU:HG	1:A:541:LEU:HD22	2.03	0.41
1:A:487:VAL:C	1:A:489:ALA:H	2.22	0.41
1:A:639:VAL:HG13	1:A:710:ILE:HD11	2.03	0.41
1:B:577:LEU:HA	1:B:577:LEU:HD23	1.88	0.41
1:A:486:ASN:HD22	1:A:487:VAL:CA	2.33	0.41
1:A:550:MET:HE2	1:A:550:MET:HB2	1.97	0.41
1:B:486:ASN:O	1:B:486:ASN:ND2	2.54	0.40
1:A:681:VAL:CG2	1:A:681:VAL:O	2.69	0.40
1:B:721:LEU:O	1:B:723:LYS:N	2.55	0.40
1:B:453:PRO:O	1:B:456:GLN:HG3	2.20	0.40
1:A:721:LEU:O	1:A:723:LYS:N	2.54	0.40
1:B:709:GLN:O	1:B:710:ILE:C	2.60	0.40
1:A:573:ILE:HG22	1:A:575:ARG:HG3	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	260/307 (85%)	235 (90%)	20 (8%)	5 (2%)	10	32
1	B	260/307 (85%)	233 (90%)	21 (8%)	6 (2%)	8	26
All	All	520/614 (85%)	468 (90%)	41 (8%)	11 (2%)	9	29

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	489	ALA
1	A	627	MET
1	B	449	ASP
1	B	489	ALA
1	B	627	MET
1	A	449	ASP
1	B	594	ASP
1	A	617	ILE
1	B	617	ILE
1	A	487	VAL
1	B	487	VAL

### 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	232/271 (86%)	210 (90%)	22 (10%)	11	30
1	B	232/271 (86%)	207 (89%)	25 (11%)	8	23
All	All	464/542 (86%)	417 (90%)	47 (10%)	9	27

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	457	ILE
1	A	485	LEU

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Mol	Chain	Res	Type
1	A	486	ASN
1	A	487	VAL
1	A	495	LEU
1	A	496	GLN
1	A	541	LEU
1	A	544	ILE
1	A	552	LYS
1	A	577	LEU
1	A	588	LEU
1	A	616	SER
1	A	618	LEU
1	A	620	MET
1	A	623	GLU
1	A	626	ARG
1	A	628	GLN
1	A	631	ASN
1	A	660	ASN
1	A	674	LEU
1	A	684	ASN
1	A	687	LYS
1	B	457	ILE
1	B	484	MET
1	B	485	LEU
1	B	486	ASN
1	B	487	VAL
1	B	495	LEU
1	B	496	GLN
1	B	541	LEU
1	B	552	LYS
1	B	577	LEU
1	B	588	LEU
1	B	616	SER
1	B	618	LEU
1	B	620	MET
1	B	623	GLU
1	B	626	ARG
1	B	628	GLN
1	B	631	ASN
1	B	658	ASN
1	B	660	ASN
1	B	662	ARG
1	B	674	LEU

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Mol	Chain	Res	Type
1	B	675	SER
1	B	684	ASN
1	B	687	LYS

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	524	GLN
1	A	628	GLN
1	A	631	ASN
1	A	660	ASN
1	A	684	ASN
1	B	461	GLN
1	B	486	ASN
1	B	524	GLN
1	B	628	GLN
1	B	631	ASN
1	B	658	ASN
1	B	660	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 ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

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	SM5	A	1	-	30,32,32	8.64	15 (50%)	35,45,45	2.16	13 (37%)
2	SM5	B	1	-	30,32,32	9.77	14 (46%)	35,45,45	1.88	11 (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	SM5	A	1	-	-	0/9/31/31	0/5/5/5
2	SM5	B	1	-	-	0/9/31/31	0/5/5/5

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	SM5	C6-C2	-15.60	1.32	1.49
2	B	1	SM5	C26-N27	-14.37	1.14	1.28
2	A	1	SM5	C23-C18	-6.49	1.35	1.52
2	A	1	SM5	C26-N27	-5.62	1.23	1.28
2	B	1	SM5	C23-C18	-4.81	1.39	1.52
2	B	1	SM5	C6-C2	-4.36	1.44	1.49
2	A	1	SM5	C1-C12	-3.70	1.43	1.49
2	B	1	SM5	C1-C2	-2.40	1.37	1.40
2	A	1	SM5	C1-C2	-2.00	1.37	1.40
2	A	1	SM5	C17-C16	2.52	1.43	1.39
2	A	1	SM5	C11-C10	2.56	1.43	1.38
2	B	1	SM5	C25-C26	2.89	1.53	1.51
2	B	1	SM5	C11-C10	3.12	1.45	1.38
2	A	1	SM5	C25-C26	3.29	1.53	1.51
2	B	1	SM5	C11-C6	3.52	1.46	1.39
2	B	1	SM5	C23-C22	3.65	1.66	1.52
2	B	1	SM5	C5-N4	3.77	1.39	1.35
2	A	1	SM5	C5-N4	3.97	1.40	1.35
2	A	1	SM5	C15-C16	4.00	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	SM5	C17-C16	4.49	1.47	1.39
2	A	1	SM5	C11-C6	4.71	1.49	1.39
2	B	1	SM5	C17-C12	6.82	1.52	1.39
2	A	1	SM5	C7-C8	7.46	1.54	1.38
2	B	1	SM5	C7-C8	8.34	1.55	1.38
2	A	1	SM5	C17-C12	8.82	1.55	1.39
2	B	1	SM5	C24-C25	10.23	1.66	1.54
2	A	1	SM5	C24-C25	11.59	1.68	1.54
2	A	1	SM5	C2-N3	39.29	1.68	1.35
2	B	1	SM5	C2-N3	47.90	1.75	1.35

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	SM5	C13-C12-C1	-5.14	113.04	120.93
2	B	1	SM5	C22-C23-C18	-3.18	107.90	110.41
2	A	1	SM5	C24-C16-C15	-3.18	109.66	111.38
2	A	1	SM5	C12-C17-C16	-3.14	117.14	121.96
2	B	1	SM5	C13-C12-C1	-2.99	116.34	120.93
2	B	1	SM5	C12-C17-C16	-2.84	117.59	121.96
2	A	1	SM5	C5-N4-C18	-2.81	122.64	125.52
2	A	1	SM5	C2-N3-N4	-2.75	102.28	105.37
2	A	1	SM5	C14-C15-C26	-2.73	122.72	132.89
2	A	1	SM5	C22-C23-C18	-2.65	108.31	110.41
2	B	1	SM5	C14-C15-C26	-2.59	123.27	132.89
2	B	1	SM5	C25-C24-C16	-2.41	102.69	104.39
2	B	1	SM5	C5-N4-C18	-2.35	123.12	125.52
2	B	1	SM5	C2-N3-N4	-2.23	102.87	105.37
2	A	1	SM5	C20-C19-C18	2.10	112.07	110.41
2	A	1	SM5	C14-C15-C16	2.22	123.55	120.64
2	A	1	SM5	C6-C2-N3	2.43	123.85	120.71
2	B	1	SM5	C17-C12-C1	2.44	124.57	120.67
2	B	1	SM5	C6-C2-N3	2.97	124.55	120.71
2	A	1	SM5	C23-C18-N4	3.17	115.33	110.80
2	B	1	SM5	C12-C1-C2	3.61	135.25	127.43
2	B	1	SM5	C23-C18-N4	3.90	116.38	110.80
2	A	1	SM5	C17-C12-C1	4.10	127.24	120.67
2	A	1	SM5	C12-C1-C2	4.51	137.20	127.43

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	SM5	6	0
2	B	1	SM5	8	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	264/307 (85%)	0.07	7 (2%) 58 45	38, 52, 89, 123	0
1	B	264/307 (85%)	0.02	11 (4%) 40 28	38, 53, 89, 123	0
All	All	528/614 (85%)	0.05	18 (3%) 49 36	38, 53, 90, 123	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	614	SER	8.5
1	A	613	LEU	6.1
1	A	722	PRO	5.9
1	B	613	LEU	5.7
1	A	723	LYS	5.6
1	A	448	ASP	5.4
1	B	723	LYS	5.0
1	B	722	PRO	5.0
1	A	614	SER	4.9
1	B	448	ASP	4.3
1	B	630	LYS	3.7
1	B	486	ASN	3.6
1	B	721	LEU	3.3
1	A	486	ASN	3.0
1	B	684	ASN	2.9
1	B	629	ASP	2.4
1	B	454	ASP	2.3
1	A	630	LYS	2.1

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

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

### 6.3 Carbohydrates [i](#)

There are no 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	SM5	A	1	28/28	0.93	0.24	1.77	41,44,49,56	0
2	SM5	B	1	28/28	0.94	0.18	0.24	39,44,46,48	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.