



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 10:32 am GMT

PDB ID : 3R7Q  
Title : Structure-based design of thienobenzoxepin inhibitors of PI3- kinase  
Authors : Murray, J.M.; Wiesmann, C.  
Deposited on : 2011-03-22  
Resolution : 2.50 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](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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

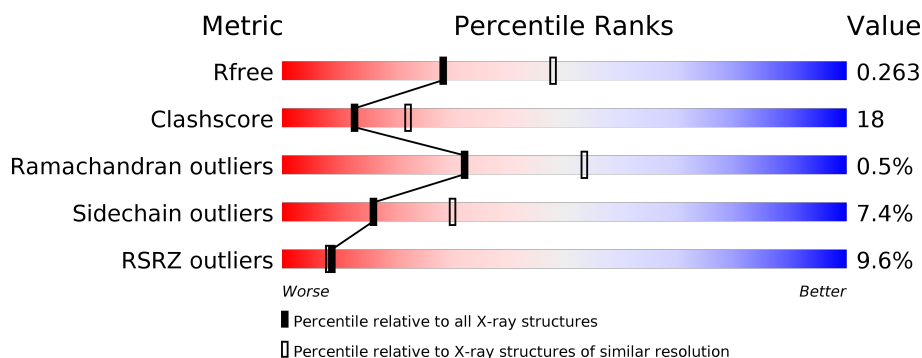
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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 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	966	<div> <div>8%</div> <div>56%</div> <div>27%</div> <div>•</div> <div>13%</div> </div>

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6800 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.

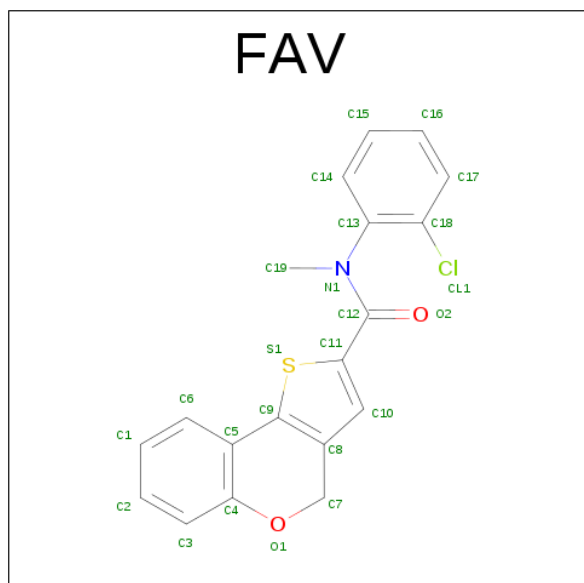
- Molecule 1 is a protein called Phosphatidylinositol-4,5-bisphosphate 3-kinase catalytic subunit gamma isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	837	Total	C	N	O	S	0	0	0
			6776	4354	1152	1234	36			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	MET	-	INITIATING METHIONINE	UNP P48736
A	1103	HIS	-	EXPRESSION TAG	UNP P48736
A	1104	HIS	-	EXPRESSION TAG	UNP P48736
A	1105	HIS	-	EXPRESSION TAG	UNP P48736
A	1106	HIS	-	EXPRESSION TAG	UNP P48736
A	1107	HIS	-	EXPRESSION TAG	UNP P48736
A	1108	HIS	-	EXPRESSION TAG	UNP P48736

- Molecule 2 is N-(2-CHLOROPHENYL)-N-METHYL-4H-THIENO[3,2-C]CHROMENE-2-CARBOXAMIDE (three-letter code: FAV) (formula: C<sub>19</sub>H<sub>14</sub>ClNO<sub>2</sub>S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	S	0	0
			24	19	1	1	2	1		



K1078	G1079	W1080	T1081	V1082	Q1083	F1084	N1085	W1086	F1087	L1088	H1089	L1090	V1091	LEU	GLY	ILE	LYS	GLN	GLY	GLU	LYS	HIS	SER	ALA	HIS	HIS	HIS	HIS	HIS
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## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	141.78Å 67.49Å 105.89Å 90.00° 96.43° 90.00°	Depositor
Resolution (Å)	19.93 – 2.50 19.93 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.5 (19.93-2.50) 98.5 (19.93-2.50)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.84 (at 2.50Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_798)	Depositor
R, $R_{free}$	0.216 , 0.269 0.209 , 0.263	Depositor DCC
$R_{free}$ test set	1709 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	66.4	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 53.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6800	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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, 0.333 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: FAV

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	0.61	0/6922	0.85	5/9363 (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	3

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	870	ILE	CG1-CB-CG2	-7.52	94.86	111.40
1	A	575	LEU	CA-CB-CG	7.47	132.47	115.30
1	A	651	LEU	CA-CB-CG	6.04	129.19	115.30
1	A	160	TYR	CB-CG-CD2	-6.03	117.38	121.00
1	A	160	TYR	CB-CG-CD1	5.27	124.16	121.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	756	LYS	Peptide
1	A	776	ASN	Peptide
1	A	777	SER	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	6776	0	6793	250	0
2	A	24	0	14	6	0
All	All	6800	0	6807	251	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 18.

All (251) 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:1035:LEU:HD12	1:A:1048:ILE:HD13	1.31	1.10
1:A:184:ARG:NH2	1:A:321:GLU:OE1	1.88	1.07
1:A:997:THR:HG23	1:A:1001:LYS:HB3	1.38	1.05
1:A:1039:MET:HB3	1:A:1040:PRO:HD2	1.40	0.99
1:A:170:ASP:OD2	1:A:476:ARG:NH2	1.98	0.96
1:A:807:LYS:HD2	1:A:807:LYS:H	1.35	0.91
1:A:1000:LYS:NZ	1:A:1073:GLU:OE1	2.07	0.87
1:A:819:ASP:OD1	1:A:821:THR:OG1	1.94	0.84
1:A:372:VAL:C	1:A:374:PRO:HD2	1.96	0.84
1:A:804:MET:HE2	1:A:831:ILE:HG12	1.57	0.84
1:A:217:ASN:HD22	1:A:219:CYS:HB2	1.43	0.84
1:A:272:LEU:HB3	1:A:305:VAL:HG11	1.63	0.79
1:A:947:ARG:NH2	1:A:963:ILE:O	2.16	0.79
1:A:741:MET:HE1	1:A:778:GLN:HB2	1.66	0.77
1:A:217:ASN:ND2	1:A:219:CYS:HB2	2.01	0.76
1:A:225:HIS:HE1	1:A:304:HIS:HD2	1.31	0.75
1:A:164:ASP:OD2	1:A:166:SER:OG	2.04	0.74
1:A:182:THR:HB	1:A:183:PRO:HD3	1.71	0.73
1:A:759:VAL:HG12	1:A:760:SER:H	1.54	0.72
1:A:564:LEU:HD11	1:A:1048:ILE:HG22	1.70	0.71
1:A:888:ILE:HD13	1:A:952:ILE:HG22	1.73	0.71
1:A:473:PHE:HD2	1:A:527:ILE:HD13	1.56	0.70
1:A:743:GLN:HG2	1:A:876:ILE:HD12	1.71	0.70
1:A:225:HIS:CE1	1:A:304:HIS:HD2	2.09	0.70
1:A:302:GLU:HB2	1:A:304:HIS:CE1	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:887:THR:HG22	1:A:890:LYS:H	1.56	0.69
1:A:759:VAL:HG12	1:A:760:SER:N	2.07	0.69
1:A:1035:LEU:HB3	1:A:1042:LEU:HG	1.74	0.69
1:A:614:ARG:HG2	1:A:617:TRP:HB3	1.75	0.69
1:A:758:ASP:OD1	1:A:759:VAL:N	2.27	0.68
1:A:202:VAL:HG12	1:A:203:THR:N	2.08	0.67
1:A:373:LEU:HD21	1:A:406:GLU:HG3	1.77	0.67
1:A:760:SER:O	1:A:763:VAL:HG12	1.95	0.67
1:A:760:SER:HB3	1:A:763:VAL:HG12	1.78	0.66
1:A:181:VAL:HG12	1:A:185:MET:CE	2.26	0.66
1:A:202:VAL:CG1	1:A:203:THR:N	2.58	0.65
1:A:743:GLN:CG	1:A:876:ILE:HD12	2.28	0.64
1:A:219:CYS:HA	1:A:235:VAL:O	1.97	0.64
1:A:905:GLU:HG2	1:A:993:PHE:CZ	2.32	0.64
1:A:1039:MET:CB	1:A:1040:PRO:HD2	2.20	0.63
1:A:214:LYS:NZ	1:A:300:GLY:HA2	2.14	0.63
1:A:807:LYS:H	1:A:807:LYS:CD	2.10	0.63
1:A:888:ILE:CD1	1:A:952:ILE:HG22	2.28	0.63
1:A:807:LYS:HD2	1:A:807:LYS:N	2.11	0.63
1:A:181:VAL:HG12	1:A:185:MET:HE2	1.81	0.63
1:A:1000:LYS:HE2	1:A:1000:LYS:HA	1.82	0.62
1:A:287:ILE:HD12	1:A:288:LYS:N	2.14	0.62
1:A:1008:LYS:O	1:A:1012:ILE:HG13	1.99	0.62
1:A:372:VAL:HB	1:A:374:PRO:HD2	1.82	0.62
1:A:886:THR:HG22	1:A:887:THR:H	1.64	0.61
1:A:888:ILE:HD13	1:A:952:ILE:CG2	2.30	0.61
1:A:879:ILE:HD12	2:A:1:FAV:CL1	2.37	0.61
1:A:758:ASP:HB3	1:A:809:LYS:HE3	1.82	0.61
1:A:746:THR:HG21	1:A:834:HIS:HB3	1.82	0.61
1:A:831:ILE:HG21	2:A:1:FAV:CL1	2.37	0.61
1:A:565:ASN:OD1	1:A:566:PRO:HD2	2.01	0.60
1:A:997:THR:HG23	1:A:1001:LYS:CB	2.24	0.60
1:A:241:PRO:O	1:A:244:ILE:HG12	2.01	0.60
1:A:244:ILE:HD11	1:A:272:LEU:CD1	2.31	0.60
1:A:251:LYS:HD3	1:A:251:LYS:O	2.01	0.59
1:A:576:TRP:CZ3	1:A:579:ARG:HD2	2.37	0.59
1:A:905:GLU:HG2	1:A:993:PHE:CE1	2.37	0.59
1:A:1056:THR:HG23	1:A:1056:THR:O	2.02	0.59
1:A:651:LEU:HD22	1:A:655:ASP:HB3	1.84	0.59
1:A:832:PHE:CZ	1:A:876:ILE:HD11	2.38	0.59
1:A:1086:TRP:CH2	1:A:1090:LEU:HD21	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:997:THR:HG22	1:A:998:SER:N	2.18	0.59
1:A:267:GLU:N	1:A:269:ASP:OD1	2.36	0.59
1:A:804:MET:HE1	2:A:1:FAV:H9	1.84	0.58
1:A:701:SER:O	1:A:705:GLN:HG2	2.03	0.58
1:A:983:VAL:HG22	1:A:984:PRO:HD2	1.84	0.58
1:A:808:LYS:HE3	1:A:836:ASP:OD1	2.04	0.57
1:A:237:PRO:HA	1:A:287:ILE:HD11	1.87	0.57
1:A:475:LEU:HD21	1:A:522:ASN:HB3	1.87	0.57
1:A:1039:MET:HB3	1:A:1040:PRO:CD	2.26	0.57
1:A:845:LEU:HD23	1:A:869:CYS:HB3	1.86	0.57
1:A:624:VAL:O	1:A:628:MET:HG2	2.05	0.57
1:A:232:THR:C	1:A:233:ILE:HD12	2.25	0.56
1:A:379:LEU:HD13	1:A:380:THR:H	1.69	0.56
1:A:225:HIS:HE1	1:A:304:HIS:CD2	2.17	0.56
1:A:759:VAL:CG1	1:A:760:SER:H	2.15	0.56
1:A:901:ALA:O	1:A:902:PHE:CD1	2.59	0.56
1:A:298:LYS:HD2	1:A:299:ASN:OD1	2.06	0.56
1:A:886:THR:HG22	1:A:887:THR:N	2.21	0.56
1:A:1006:PHE:O	1:A:1009:PHE:HB3	2.06	0.55
1:A:286:PRO:O	1:A:289:ASN:HB2	2.05	0.55
1:A:804:MET:HE3	1:A:810:PRO:HG2	1.88	0.55
1:A:746:THR:O	1:A:750:LYS:HB2	2.06	0.55
1:A:804:MET:CE	1:A:831:ILE:HG12	2.32	0.55
1:A:271:VAL:HG23	1:A:282:VAL:HG12	1.88	0.55
1:A:240:THR:HG23	1:A:241:PRO:HD2	1.88	0.55
1:A:225:HIS:CE1	1:A:304:HIS:CD2	2.95	0.55
1:A:373:LEU:N	1:A:374:PRO:HD2	2.22	0.55
1:A:921:PHE:O	1:A:925:VAL:HG23	2.06	0.54
1:A:853:SER:O	1:A:857:THR:HG23	2.07	0.54
1:A:389:HIS:O	1:A:392:GLN:HB3	2.08	0.54
1:A:360:LYS:HE2	1:A:417:SER:HA	1.89	0.54
1:A:903:LYS:O	1:A:906:VAL:HG23	2.08	0.53
1:A:1042:LEU:O	1:A:1042:LEU:HD13	2.08	0.53
1:A:1032:SER:HB3	1:A:1048:ILE:CG2	2.38	0.53
1:A:888:ILE:HD12	1:A:952:ILE:O	2.09	0.53
1:A:887:THR:HG22	1:A:889:ALA:N	2.23	0.53
1:A:273:ARG:NH2	1:A:821:THR:OG1	2.42	0.53
1:A:146:GLU:HB3	1:A:319:ARG:HH22	1.74	0.52
1:A:363:VAL:HG23	1:A:520:LEU:CD1	2.38	0.52
1:A:964:ASP:CB	2:A:1:FAV:H12	2.38	0.52
1:A:903:LYS:HB2	1:A:906:VAL:CG2	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:THR:O	1:A:230:SER:HB3	2.10	0.52
1:A:583:LEU:O	1:A:583:LEU:HD23	2.10	0.52
1:A:983:VAL:CG2	1:A:984:PRO:HD2	2.39	0.52
1:A:1043:THR:N	1:A:1047:ASP:OD2	2.41	0.52
1:A:576:TRP:CH2	1:A:579:ARG:HD2	2.45	0.52
1:A:741:MET:CE	1:A:778:GLN:HB2	2.37	0.52
1:A:144:SER:OG	1:A:146:GLU:HB2	2.09	0.51
1:A:220:ILE:O	1:A:235:VAL:N	2.43	0.51
1:A:497:PHE:HB2	1:A:1041:GLN:HG3	1.92	0.51
1:A:758:ASP:HB3	1:A:809:LYS:CE	2.40	0.51
1:A:395:CYS:HB3	1:A:416:PHE:HD2	1.75	0.51
1:A:875:LYS:N	1:A:875:LYS:HD3	2.25	0.51
2:A:1:FAV:C13	2:A:1:FAV:S1	2.98	0.51
1:A:213:LYS:HD3	1:A:214:LYS:N	2.25	0.51
1:A:1042:LEU:O	1:A:1042:LEU:HD22	2.11	0.51
1:A:752:LEU:O	1:A:753:SER:HB2	2.10	0.51
1:A:777:SER:CB	1:A:778:GLN:HA	2.40	0.51
1:A:233:ILE:N	1:A:233:ILE:HD12	2.26	0.51
1:A:311:PRO:O	1:A:313:PRO:HD3	2.12	0.50
1:A:1035:LEU:HA	1:A:1039:MET:HG2	1.94	0.50
1:A:912:LYS:NZ	1:A:918:GLU:OE2	2.41	0.50
1:A:887:THR:CG2	1:A:889:ALA:H	2.25	0.50
1:A:759:VAL:CG1	1:A:760:SER:N	2.72	0.50
1:A:902:PHE:HD2	1:A:1080:TRP:HB3	1.77	0.50
1:A:1042:LEU:CD1	1:A:1042:LEU:H	2.26	0.49
1:A:804:MET:CE	1:A:831:ILE:HG23	2.43	0.49
1:A:1002:THR:O	1:A:1003:SER:HB3	2.13	0.49
1:A:774:LEU:O	1:A:776:ASN:N	2.46	0.49
1:A:583:LEU:CD1	1:A:610:LEU:HD22	2.42	0.49
1:A:244:ILE:HD11	1:A:272:LEU:HD11	1.94	0.49
1:A:692:GLY:HA3	1:A:720:TYR:OH	2.13	0.48
1:A:302:GLU:HB2	1:A:304:HIS:HE1	1.76	0.48
1:A:804:MET:HE2	1:A:831:ILE:HG23	1.94	0.48
1:A:848:LEU:HA	1:A:848:LEU:HD12	1.63	0.48
1:A:1039:MET:CB	1:A:1040:PRO:CD	2.90	0.48
1:A:1032:SER:HB3	1:A:1048:ILE:HG21	1.95	0.48
1:A:285:THR:HG22	1:A:289:ASN:HB2	1.96	0.48
1:A:270:PHE:HB3	1:A:307:LEU:HD11	1.96	0.48
1:A:221:PHE:HB3	1:A:232:THR:HG22	1.96	0.48
1:A:887:THR:HG22	1:A:889:ALA:H	1.79	0.48
1:A:1017:TYR:CE2	1:A:1021:ARG:HD2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:PHE:HE2	1:A:234:LYS:HG2	1.79	0.47
1:A:1048:ILE:O	1:A:1051:ILE:HG22	2.14	0.47
1:A:293:VAL:O	1:A:297:LEU:HG	2.15	0.47
1:A:842:MET:HG2	1:A:869:CYS:O	2.14	0.47
1:A:472:ARG:O	1:A:473:PHE:HB2	2.14	0.47
1:A:753:SER:OG	1:A:754:ALA:N	2.46	0.47
1:A:767:LEU:HD22	1:A:771:LEU:HG	1.97	0.47
1:A:180:LEU:C	1:A:183:PRO:HD2	2.35	0.47
1:A:463:TYR:CE1	1:A:501:LYS:HA	2.50	0.47
1:A:175:PHE:HA	1:A:178:ARG:NH1	2.30	0.46
1:A:988:THR:HB	1:A:989:PRO:HD2	1.96	0.46
1:A:214:LYS:NZ	1:A:297:LEU:HA	2.30	0.46
1:A:372:VAL:HB	1:A:374:PRO:CD	2.44	0.46
1:A:379:LEU:HD13	1:A:380:THR:HG22	1.98	0.46
1:A:1032:SER:HA	1:A:1048:ILE:HD12	1.98	0.46
1:A:202:VAL:CG1	1:A:203:THR:H	2.28	0.46
1:A:214:LYS:HZ1	1:A:300:GLY:HA2	1.79	0.46
1:A:867:TYR:CE2	1:A:963:ILE:HG22	2.51	0.46
1:A:760:SER:C	1:A:762:GLN:N	2.68	0.46
1:A:1026:LEU:O	1:A:1029:ILE:HG22	2.15	0.45
1:A:1002:THR:HG22	1:A:1003:SER:H	1.81	0.45
1:A:925:VAL:O	1:A:929:VAL:HG23	2.16	0.45
1:A:721:LEU:HA	1:A:721:LEU:HD23	1.67	0.45
1:A:832:PHE:CE1	1:A:876:ILE:HD11	2.51	0.45
1:A:210:TYR:CE1	1:A:211:LEU:HG	2.51	0.45
1:A:876:ILE:HG21	1:A:876:ILE:HD13	1.64	0.45
1:A:207:LEU:HD23	1:A:208:PRO:HD2	1.99	0.45
1:A:210:TYR:O	1:A:213:LYS:HD2	2.17	0.45
1:A:220:ILE:N	1:A:235:VAL:O	2.48	0.45
1:A:660:LEU:O	1:A:664:VAL:HG23	2.16	0.45
1:A:929:VAL:HG22	1:A:995:MET:HG2	1.99	0.45
1:A:569:ALA:O	1:A:573:GLU:HG3	2.17	0.45
1:A:770:LYS:O	1:A:774:LEU:HG	2.17	0.45
1:A:213:LYS:NZ	1:A:214:LYS:HB2	2.32	0.45
1:A:236:SER:O	1:A:238:ASP:N	2.50	0.45
1:A:760:SER:HB3	1:A:763:VAL:CG1	2.47	0.45
1:A:804:MET:HB2	1:A:810:PRO:HG2	1.99	0.45
1:A:872:THR:OG1	1:A:877:GLY:HA2	2.17	0.45
1:A:892:GLN:HE21	1:A:902:PHE:HD1	1.64	0.45
1:A:1018:LEU:HD23	1:A:1021:ARG:HD3	1.98	0.44
1:A:287:ILE:HG13	1:A:287:ILE:H	1.67	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:CYS:N	1:A:524:CYS:HG	2.16	0.44
1:A:935:TYR:CE1	1:A:961:PHE:HA	2.53	0.44
1:A:904:ASP:O	1:A:990:ASP:HA	2.18	0.44
1:A:1067:TYR:O	1:A:1071:GLN:HG2	2.17	0.44
1:A:213:LYS:HD3	1:A:214:LYS:H	1.83	0.44
1:A:372:VAL:CB	1:A:374:PRO:HD2	2.45	0.44
1:A:737:GLN:O	1:A:741:MET:HG3	2.17	0.44
1:A:1045:LYS:O	1:A:1049:GLU:OE1	2.35	0.44
1:A:651:LEU:HD22	1:A:655:ASP:CB	2.47	0.44
1:A:198:MET:SD	1:A:282:VAL:HG11	2.58	0.44
1:A:304:HIS:CG	1:A:823:LEU:HD21	2.53	0.43
1:A:997:THR:CG2	1:A:998:SER:N	2.81	0.43
1:A:990:ASP:OD1	1:A:990:ASP:N	2.51	0.43
1:A:1055:LEU:O	1:A:1056:THR:HG22	2.19	0.43
1:A:739:ILE:HG13	1:A:740:GLU:N	2.33	0.43
1:A:315:LEU:O	1:A:727:ALA:HB2	2.19	0.43
1:A:915:SER:HA	1:A:916:PRO:HD3	1.89	0.43
1:A:211:LEU:HD23	1:A:211:LEU:HA	1.87	0.43
1:A:784:ARG:HG2	1:A:784:ARG:NH1	2.34	0.43
1:A:1042:LEU:HD22	1:A:1042:LEU:C	2.39	0.43
1:A:1082:VAL:HA	1:A:1085:ASN:ND2	2.33	0.43
1:A:992:LEU:HA	1:A:992:LEU:HD23	1.88	0.43
1:A:888:ILE:HG22	1:A:949:ASN:OD1	2.18	0.43
1:A:309:THR:HA	1:A:310:PRO:HD3	1.90	0.42
1:A:675:SER:O	1:A:679:ARG:HG3	2.19	0.42
1:A:777:SER:HB2	1:A:778:GLN:HA	2.00	0.42
1:A:221:PHE:HA	1:A:233:ILE:O	2.19	0.42
1:A:373:LEU:N	1:A:374:PRO:CD	2.83	0.42
1:A:390:GLY:N	1:A:636:SER:OG	2.45	0.42
1:A:774:LEU:C	1:A:776:ASN:N	2.72	0.42
1:A:246:GLN:C	1:A:248:PHE:H	2.21	0.42
1:A:590:PRO:HD3	1:A:626:LEU:HD21	2.02	0.42
1:A:498:ASN:OD1	1:A:498:ASN:C	2.57	0.42
1:A:1068:PHE:O	1:A:1071:GLN:HB2	2.20	0.42
1:A:270:PHE:HB3	1:A:307:LEU:CD1	2.49	0.42
1:A:235:VAL:HG12	1:A:236:SER:N	2.35	0.42
1:A:467:LEU:O	1:A:476:ARG:HD2	2.20	0.42
1:A:583:LEU:HD12	1:A:610:LEU:HD22	2.01	0.42
1:A:662:GLN:HE22	1:A:850:ILE:CD1	2.33	0.42
1:A:862:LEU:CD2	1:A:862:LEU:N	2.83	0.42
1:A:207:LEU:CD2	1:A:208:PRO:HD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:379:LEU:CD1	1:A:380:THR:H	2.32	0.41
1:A:359:ARG:HD3	1:A:359:ARG:HA	1.92	0.41
1:A:702:GLU:O	1:A:706:SER:HB3	2.21	0.41
1:A:887:THR:CG2	1:A:889:ALA:HB3	2.50	0.41
1:A:224:ILE:HA	1:A:305:VAL:O	2.20	0.41
1:A:942:LEU:HD23	1:A:942:LEU:HA	1.84	0.41
1:A:214:LYS:O	1:A:214:LYS:HG2	2.21	0.41
1:A:476:ARG:HB3	1:A:520:LEU:HD23	2.03	0.41
1:A:249:PHE:CD1	1:A:268:GLN:OE1	2.74	0.41
1:A:1032:SER:HB3	1:A:1048:ILE:HG23	2.03	0.41
1:A:373:LEU:CD2	1:A:406:GLU:HG3	2.49	0.41
1:A:808:LYS:O	1:A:810:PRO:HD3	2.21	0.41
1:A:245:LEU:HA	1:A:245:LEU:HD23	1.91	0.41
1:A:287:ILE:HA	1:A:290:PHE:HD1	1.86	0.41
1:A:227:SER:C	1:A:229:THR:H	2.24	0.40
1:A:169:HIS:CE1	1:A:466:LEU:CD1	3.04	0.40
1:A:964:ASP:CG	2:A:1:FAV:H12	2.41	0.40
1:A:304:HIS:CE1	1:A:823:LEU:HD21	2.57	0.40
1:A:182:THR:CB	1:A:183:PRO:HD3	2.45	0.40
1:A:840:GLN:O	1:A:844:ILE:HG12	2.21	0.40
1:A:184:ARG:HD3	1:A:719:ALA:O	2.22	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	819/966 (85%)	775 (95%)	40 (5%)	4 (0%)	32 53

All (4) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	775	GLN
1	A	759	VAL
1	A	1040	PRO
1	A	237	PRO

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	748/864 (87%)	693 (93%)	55 (7%)	16	30

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	GLU
1	A	146	GLU
1	A	147	SER
1	A	207	LEU
1	A	213	LYS
1	A	227	SER
1	A	228	THR
1	A	238	ASP
1	A	252	MET
1	A	269	ASP
1	A	298	LYS
1	A	373	LEU
1	A	379	LEU
1	A	381	VAL
1	A	435	CYS
1	A	476	ARG
1	A	477	ARG
1	A	511	GLU
1	A	520	LEU
1	A	521	ASP
1	A	523	TYR
1	A	549	ASN
1	A	550	GLN

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Mol	Chain	Res	Type
1	A	568	THR
1	A	570	GLU
1	A	575	LEU
1	A	601	GLN
1	A	610	LEU
1	A	626	LEU
1	A	647	LYS
1	A	682	LEU
1	A	707	ARG
1	A	717	LEU
1	A	739	ILE
1	A	760	SER
1	A	765	SER
1	A	767	LEU
1	A	778	GLN
1	A	838	LEU
1	A	843	LEU
1	A	845	LEU
1	A	848	LEU
1	A	865	LEU
1	A	887	THR
1	A	890	LYS
1	A	907	LEU
1	A	926	GLU
1	A	959	ASN
1	A	967	HIS
1	A	1026	LEU
1	A	1042	LEU
1	A	1051	ILE
1	A	1078	LYS
1	A	1086	TRP
1	A	1088	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	225	HIS
1	A	304	HIS
1	A	778	GLN
1	A	1085	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 ⓘ

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$
2	FAV	A	1	-	20,27,27	0.69	1 (5%)	28,39,39	1.68	5 (17%)

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	FAV	A	1	-	-	0/8/21/21	0/4/4/4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	FAV	C13-N1	-2.34	1.41	1.44

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	FAV	C4-C5-C9	-2.50	115.96	118.56
2	A	1	FAV	C11-C12-N1	2.45	124.33	118.50
2	A	1	FAV	O1-C4-C3	2.93	121.58	116.76
2	A	1	FAV	C6-C5-C4	3.12	121.72	117.38
2	A	1	FAV	C7-O1-C4	4.74	124.67	113.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	FAV	6	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

### 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<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	837/966 (86%)	0.39	80 (9%) <b>9</b> <b>8</b>	43, 85, 155, 227	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	528	ALA	16.6
1	A	1089	HIS	11.5
1	A	322	GLU	9.9
1	A	1086	TRP	8.6
1	A	756	LYS	8.4
1	A	544	ARG	8.1
1	A	143	MET	7.8
1	A	1000	LYS	7.5
1	A	1045	LYS	7.0
1	A	374	PRO	6.8
1	A	1044	SER	6.5
1	A	378	ASP	6.3
1	A	377	THR	6.2
1	A	1091	VAL	6.2
1	A	757	TYR	6.2
1	A	253	ALA	5.7
1	A	216	ALA	5.6
1	A	1046	GLU	5.4
1	A	148	GLN	4.7
1	A	758	ASP	4.7
1	A	981	GLU	4.6
1	A	999	GLY	4.5
1	A	1043	THR	4.4
1	A	982	ARG	4.3
1	A	250	THR	4.3
1	A	523	TYR	4.1
1	A	759	VAL	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	267	GLU	4.0
1	A	823	LEU	4.0
1	A	987	LEU	3.7
1	A	998	SER	3.5
1	A	919	GLU	3.5
1	A	1090	LEU	3.4
1	A	1062	GLU	3.4
1	A	212	TRP	3.4
1	A	249	PHE	3.3
1	A	1084	PHE	3.3
1	A	967	HIS	3.2
1	A	373	LEU	3.2
1	A	916	PRO	3.2
1	A	1041	GLN	3.1
1	A	991	PHE	3.1
1	A	488	SER	3.0
1	A	489	GLY	2.9
1	A	215	ILE	2.9
1	A	269	ASP	2.8
1	A	270	PHE	2.8
1	A	932	CYS	2.8
1	A	909	HIS	2.8
1	A	1007	GLN	2.7
1	A	252	MET	2.7
1	A	210	TYR	2.7
1	A	201	TRP	2.7
1	A	1040	PRO	2.7
1	A	809	LYS	2.6
1	A	1015	LYS	2.5
1	A	1088	LEU	2.5
1	A	859	SER	2.5
1	A	1042	LEU	2.4
1	A	930	TYR	2.4
1	A	778	GLN	2.4
1	A	1082	VAL	2.4
1	A	777	SER	2.4
1	A	826	GLU	2.4
1	A	751	SER	2.4
1	A	832	PHE	2.4
1	A	217	ASN	2.3
1	A	808	LYS	2.2
1	A	315	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	431	LEU	2.2
1	A	776	ASN	2.2
1	A	1087	PHE	2.2
1	A	768	LYS	2.1
1	A	527	ILE	2.1
1	A	522	ASN	2.1
1	A	825	ASN	2.1
1	A	1004	PRO	2.1
1	A	755	GLU	2.1
1	A	268	GLN	2.0
1	A	613	ARG	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 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(Å <sup>2</sup> )	Q < 0.9
2	FAV	A	1	24/24	0.97	0.11	-0.85	65,78,88,294	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.