



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:54 AM GMT

PDB ID : 2ETM
Title : Crystal Structure of Focal Adhesion Kinase Domain Complexed with 7H-Pyrrolo [2,3-d] pyrimidine Derivative
Authors : Lee, C.C.
Deposited on : 2005-10-27
Resolution : 2.30 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

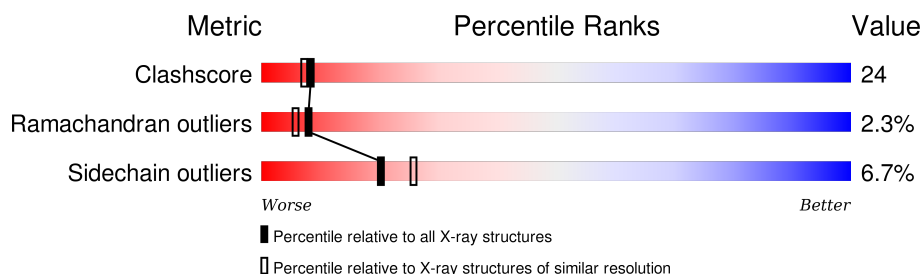
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	281	
1	B	281	

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	7PY	B	132	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4400 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.

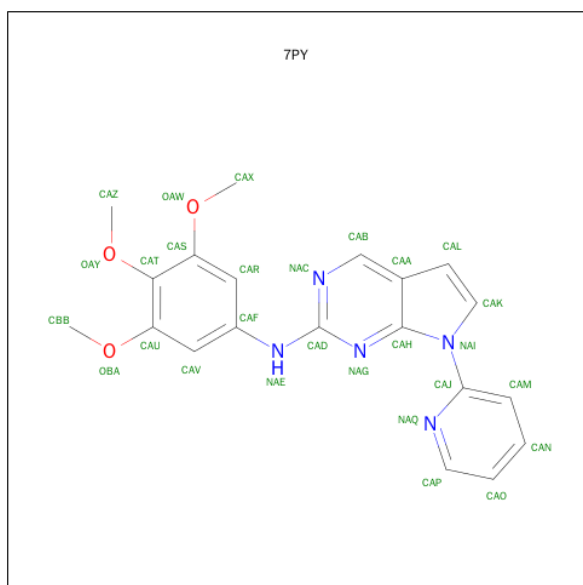
- Molecule 1 is a protein called Focal adhesion kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	262	Total	C	N	O	S	0	0	0
			2111	1349	365	379	18			
1	B	263	Total	C	N	O	S	0	0	0
			2114	1346	366	385	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	409	GLY	-	CLONING ARTIFACT	UNP Q05397
A	410	ALA	-	CLONING ARTIFACT	UNP Q05397
B	409	GLY	-	CLONING ARTIFACT	UNP Q05397
B	410	ALA	-	CLONING ARTIFACT	UNP Q05397

- Molecule 2 is 7-PYRIDIN-2-YL-N-(3,4,5-TRIMETHOXYPHENYL)-7H-PYRROLO[2,3-D]PYRIMIDIN-2-AMINE (three-letter code: 7PY) (formula: C₂₀H₁₉N₅O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			28	20	5	3		
2	B	1	Total	C	N	O	0	0
			28	20	5	3		

- Molecule 3 is water.

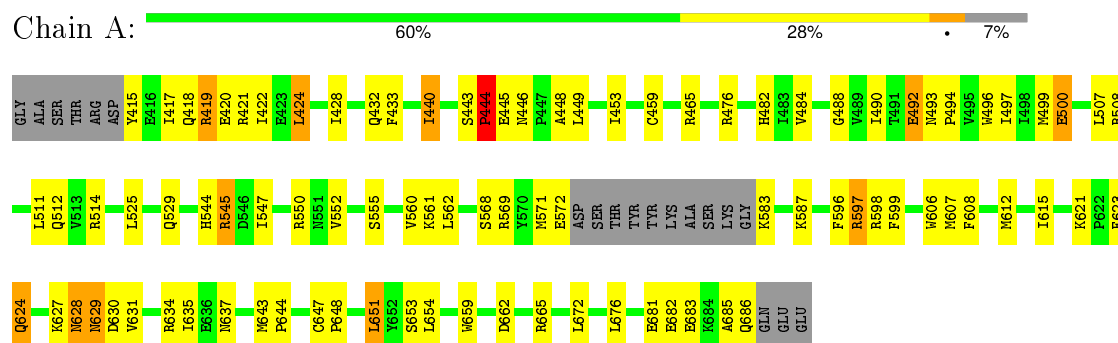
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	70	Total	O	0	0
			70	70		
3	B	49	Total	O	0	0
			49	49		

3 Residue-property plots

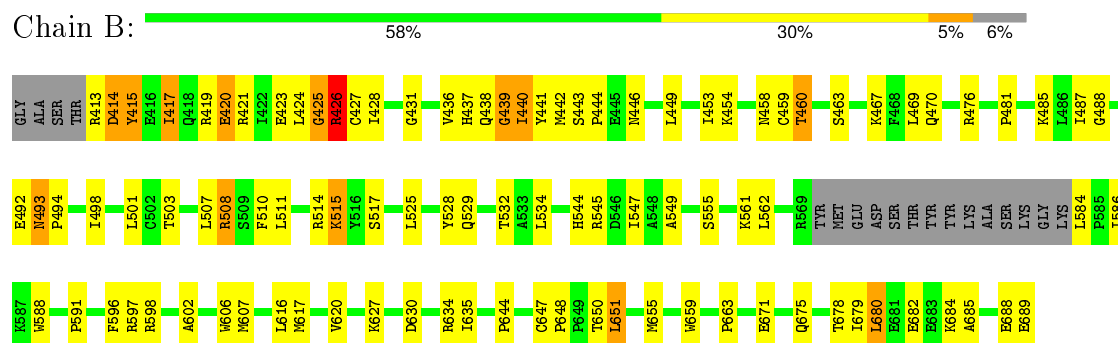
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

- Molecule 1: Focal adhesion kinase 1



- Molecule 1: Focal adhesion kinase 1



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	45.70Å 51.92Å 66.88Å 100.08° 103.94° 90.17°	Depositor
Resolution (Å)	6.00 – 2.30	Depositor
% Data completeness (in resolution range)	97.2 (6.00-2.30)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	0.07	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.210 , 0.295	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4400	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 7PY

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	0.47	0/2158	0.71	0/2916
1	B	0.46	0/2160	0.71	0/2919
All	All	0.46	0/4318	0.71	0/5835

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2111	0	2123	104	0
1	B	2114	0	2112	93	0
2	A	28	0	19	7	0
2	B	28	0	19	9	0
3	A	70	0	0	5	0
3	B	49	0	0	6	0
All	All	4400	0	4273	207	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 (207) 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:587:LYS:HD3	1:A:623:PHE:HB2	1.44	0.97
1:A:621:LYS:HD2	1:A:624:GLN:HE22	1.32	0.95
1:A:421:ARG:HH12	1:A:444:PRO:HD2	1.31	0.95
1:A:547:ILE:H	1:A:607:MET:HE3	1.37	0.90
1:B:616:LEU:HD11	1:B:651:LEU:HD13	1.55	0.88
1:A:621:LYS:HD2	1:A:624:GLN:NE2	1.91	0.84
1:A:417:ILE:HG22	1:A:418:GLN:H	1.43	0.84
1:A:648:PRO:HG2	1:A:651:LEU:HB2	1.59	0.84
2:B:132:7PY:HAZ2	3:B:55:HOH:O	1.77	0.84
1:B:525:LEU:O	1:B:529:GLN:HG3	1.77	0.83
1:B:428:ILE:O	2:B:132:7PY:HAX2	1.79	0.82
1:A:424:LEU:HD12	1:A:424:LEU:H	1.45	0.81
1:A:432:GLN:HE22	1:A:571:MET:HG3	1.45	0.80
1:A:421:ARG:NH1	1:A:443:SER:HA	1.95	0.79
1:A:432:GLN:NE2	1:A:571:MET:HG3	1.98	0.79
1:A:612:MET:HE1	1:A:676:LEU:HD11	1.66	0.77
1:B:419:ARG:HH11	1:B:419:ARG:HG2	1.49	0.77
1:A:587:LYS:HE2	1:A:623:PHE:O	1.84	0.76
2:B:132:7PY:HAZ3	2:B:132:7PY:HBB2	1.69	0.75
1:A:525:LEU:O	1:A:529:GLN:HG3	1.85	0.75
1:A:421:ARG:HH12	1:A:444:PRO:CD	1.99	0.75
1:B:413:ARG:O	1:B:415:TYR:N	2.20	0.74
2:A:131:7PY:HAZ3	2:A:131:7PY:HBB2	1.69	0.74
2:A:131:7PY:NAG	2:A:131:7PY:HAR	2.03	0.73
1:B:684:LYS:O	1:B:688:GLU:HG3	1.88	0.73
1:A:630:ASP:O	1:A:634:ARG:HG3	1.89	0.72
1:A:422:ILE:HD12	1:A:490:ILE:HD11	1.73	0.70
1:B:425:GLY:O	1:B:426:ARG:HB3	1.90	0.69
1:A:552:VAL:CG1	1:A:560:VAL:HG13	2.21	0.69
1:B:438:GLN:O	1:B:439:GLY:O	2.11	0.69
1:A:417:ILE:HG22	1:A:418:GLN:N	2.09	0.66
2:B:132:7PY:NAG	2:B:132:7PY:HAR	2.11	0.66
1:A:662:ASP:H	1:A:665:ARG:NH2	1.94	0.66
1:B:547:ILE:H	1:B:607:MET:HE3	1.60	0.66
1:A:612:MET:CE	1:A:676:LEU:HD11	2.25	0.65
1:B:671:GLU:HG2	1:B:675:GLN:HE21	1.61	0.65
1:B:544:HIS:O	1:B:545:ARG:HB2	1.95	0.65
1:B:443:SER:HB2	1:B:446:ASN:O	1.97	0.65
1:B:644:PRO:HG2	1:B:647:CYS:HB2	1.78	0.65
1:B:436:VAL:HG22	1:B:454:LYS:HG2	1.78	0.64
1:A:433:PHE:HZ	1:A:571:MET:HE3	1.61	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:584:LEU:O	1:B:586:ILE:HD12	1.98	0.64
1:B:493:ASN:N	1:B:493:ASN:HD22	1.95	0.63
1:A:571:MET:O	1:A:572:GLU:HB2	1.98	0.63
1:A:418:GLN:NE2	1:A:421:ARG:HE	1.95	0.63
1:B:436:VAL:HG21	2:B:132:7PY:CAP	2.29	0.63
1:A:662:ASP:HB3	1:A:665:ARG:NE	2.14	0.63
1:A:432:GLN:OE1	1:A:571:MET:HG3	1.99	0.63
1:A:683:GLU:OE1	1:A:683:GLU:HA	1.99	0.62
1:B:421:ARG:CZ	1:B:444:PRO:HG3	2.29	0.62
1:A:685:ALA:O	1:A:686:GLN:HG3	1.98	0.62
1:A:612:MET:HE2	1:A:676:LEU:HD21	1.81	0.62
1:B:586:ILE:H	1:B:586:ILE:HD12	1.65	0.62
1:A:421:ARG:NH1	1:A:444:PRO:HD2	2.11	0.61
1:B:678:THR:O	1:B:682:GLU:HG3	2.00	0.61
1:A:497:ILE:HG22	1:A:499:MET:HE3	1.82	0.61
1:B:428:ILE:O	2:B:132:7PY:CAX	2.49	0.61
1:B:481:PRO:O	1:B:561:LYS:HE2	2.01	0.61
1:A:421:ARG:HH11	1:A:443:SER:HA	1.66	0.61
1:A:419:ARG:NH1	1:A:494:PRO:HD2	2.16	0.61
1:A:440:ILE:HD13	1:A:448:ALA:HB1	1.82	0.60
1:A:432:GLN:HE22	1:A:571:MET:CG	2.13	0.60
1:A:419:ARG:HH12	1:A:493:ASN:C	2.04	0.60
1:A:421:ARG:NH1	1:A:444:PRO:CD	2.64	0.60
1:B:528:TYR:O	1:B:532:THR:HG23	2.02	0.60
1:A:428:ILE:O	2:A:131:7PY:HAX2	2.02	0.60
1:A:444:PRO:C	1:A:446:ASN:H	2.05	0.59
1:B:413:ARG:C	1:B:415:TYR:H	2.05	0.59
1:A:552:VAL:HG13	1:A:561:LYS:O	2.02	0.59
1:A:547:ILE:N	1:A:607:MET:HE3	2.13	0.59
1:B:616:LEU:HD13	1:B:647:CYS:SG	2.42	0.59
1:A:432:GLN:CD	1:A:571:MET:HG3	2.23	0.59
2:B:132:7PY:HAN	2:B:132:7PY:HAX3	1.85	0.58
1:B:596:PHE:HB2	1:B:598:ARG:HD3	1.83	0.58
1:A:508:ARG:O	1:A:512:GLN:HG3	2.04	0.58
1:A:415:TYR:CE1	1:A:476:ARG:HG3	2.39	0.58
1:A:545:ARG:C	1:A:607:MET:HE1	2.24	0.57
1:B:648:PRO:HG2	1:B:651:LEU:HB2	1.86	0.57
1:B:510:PHE:O	1:B:514:ARG:HG2	2.04	0.57
1:B:492:GLU:OE2	1:B:492:GLU:HA	2.04	0.57
1:A:544:HIS:O	1:A:545:ARG:HB2	2.05	0.56
1:B:515:LYS:HB3	1:B:515:LYS:NZ	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:654:LEU:HD11	1:A:672:LEU:HD23	1.88	0.55
1:A:624:GLN:OE1	1:A:624:GLN:HA	2.07	0.55
1:B:440:ILE:HG13	1:B:441:TYR:N	2.21	0.54
1:B:650:THR:HG22	1:B:679:ILE:HD11	1.90	0.54
1:A:432:GLN:NE2	1:A:568:SER:O	2.41	0.54
2:A:131:7PY:NAG	2:A:131:7PY:CAR	2.66	0.54
1:A:631:VAL:O	1:A:635:ILE:HG12	2.07	0.54
1:A:492:GLU:O	1:A:493:ASN:HB2	2.07	0.53
1:A:444:PRO:O	1:A:446:ASN:N	2.42	0.53
1:A:552:VAL:HG11	1:A:560:VAL:HG13	1.91	0.52
1:B:443:SER:CB	1:B:446:ASN:HB2	2.39	0.52
1:A:545:ARG:HD3	1:A:599:PHE:CG	2.45	0.52
1:A:418:GLN:CG	1:A:421:ARG:HD2	2.40	0.52
1:B:617:MET:O	1:B:620:VAL:HG23	2.09	0.52
1:B:453:ILE:HG12	1:B:498:ILE:HD12	1.91	0.51
1:A:418:GLN:NE2	1:A:421:ARG:NE	2.59	0.51
1:A:608:PHE:CZ	1:A:612:MET:HE3	2.45	0.51
1:A:612:MET:CE	1:A:676:LEU:HD21	2.39	0.51
1:B:415:TYR:CE1	1:B:476:ARG:HG3	2.46	0.51
1:B:417:ILE:HD11	1:B:488:GLY:HA3	1.91	0.51
1:B:421:ARG:HD3	1:B:442:MET:O	2.11	0.51
1:A:628:ASN:O	1:A:631:VAL:N	2.44	0.50
1:B:663:PRO:HB3	3:B:117:HOH:O	2.10	0.50
1:B:425:GLY:O	1:B:426:ARG:CB	2.57	0.50
1:B:616:LEU:CD1	1:B:655:MET:HE1	2.41	0.50
1:A:550:ARG:HA	3:A:109:HOH:O	2.11	0.50
1:B:617:MET:SD	1:B:644:PRO:HB3	2.52	0.49
1:B:420:GLU:H	1:B:420:GLU:CD	2.15	0.49
1:A:545:ARG:HD3	1:A:599:PHE:CD2	2.47	0.49
1:B:476:ARG:HG2	1:B:476:ARG:O	2.12	0.49
1:A:493:ASN:HA	1:A:494:PRO:C	2.33	0.49
1:A:428:ILE:HD12	2:A:131:7PY:CAU	2.43	0.49
1:B:424:LEU:H	1:B:424:LEU:HD23	1.76	0.49
1:A:628:ASN:O	1:A:630:ASP:N	2.46	0.49
1:A:514:ARG:HG3	1:A:514:ARG:HH11	1.77	0.48
1:A:453:ILE:N	1:A:453:ILE:HD12	2.28	0.48
1:A:508:ARG:HB2	3:A:109:HOH:O	2.14	0.48
1:B:508:ARG:HD2	1:B:549:ALA:HB3	1.95	0.48
1:A:428:ILE:CG2	1:A:428:ILE:O	2.61	0.48
1:B:588:TRP:O	1:B:607:MET:HG2	2.14	0.48
1:B:508:ARG:HH11	1:B:508:ARG:HG3	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:497:ILE:CG2	1:A:499:MET:HE3	2.43	0.48
1:B:606:TRP:CE3	1:B:659:TRP:HA	2.47	0.48
1:B:427:CYS:HB2	1:B:437:HIS:CE1	2.49	0.47
1:A:596:PHE:C	1:A:597:ARG:HD2	2.34	0.47
1:B:680:LEU:O	1:B:684:LYS:HG3	2.14	0.47
1:B:426:ARG:HE	1:B:438:GLN:HE22	1.62	0.47
1:A:421:ARG:HH12	1:A:443:SER:HA	1.75	0.47
1:B:616:LEU:CD1	1:B:651:LEU:HD13	2.36	0.47
1:B:419:ARG:NH1	1:B:419:ARG:HG2	2.26	0.47
1:B:414:ASP:OD2	1:B:476:ARG:NH2	2.47	0.47
1:A:662:ASP:H	1:A:665:ARG:HH21	1.62	0.47
1:B:547:ILE:N	1:B:607:MET:HE3	2.27	0.47
1:B:421:ARG:NE	1:B:444:PRO:HG3	2.29	0.47
1:A:682:GLU:O	1:A:686:GLN:HB2	2.15	0.47
1:B:634:ARG:HG3	1:B:634:ARG:HH11	1.80	0.47
1:B:426:ARG:NE	1:B:438:GLN:HE22	2.13	0.46
1:B:515:LYS:HB3	1:B:515:LYS:HZ2	1.80	0.46
1:A:490:ILE:HD12	1:A:496:TRP:CE3	2.51	0.46
1:B:463:SER:O	1:B:467:LYS:HG3	2.14	0.46
2:B:132:7PY:NAG	2:B:132:7PY:CAR	2.70	0.46
1:A:415:TYR:HB3	1:A:488:GLY:HA2	1.96	0.46
1:B:501:LEU:HD13	1:B:503:THR:N	2.31	0.46
1:B:685:ALA:O	1:B:689:GLU:HG3	2.16	0.46
1:B:431:GLY:HA3	3:B:116:HOH:O	2.15	0.45
1:A:417:ILE:CG2	1:A:418:GLN:N	2.80	0.45
1:A:417:ILE:CG2	1:A:418:GLN:H	2.21	0.45
1:B:458:ASN:C	1:B:460:THR:H	2.20	0.45
1:A:569:ARG:O	1:A:583:LYS:O	2.35	0.45
1:A:484:VAL:HG23	1:A:562:LEU:O	2.17	0.45
1:B:507:LEU:O	1:B:511:LEU:HG	2.17	0.44
1:A:424:LEU:CD1	1:A:424:LEU:H	2.24	0.44
1:B:591:PRO:HG3	1:B:635:ILE:HG21	2.00	0.44
1:B:534:LEU:HA	1:B:534:LEU:HD23	1.67	0.44
1:A:627:LYS:O	1:A:630:ASP:HB2	2.17	0.44
1:B:421:ARG:CD	1:B:444:PRO:HG3	2.48	0.44
1:A:507:LEU:HD21	1:A:615:ILE:HG12	1.99	0.44
1:B:493:ASN:HA	1:B:494:PRO:C	2.38	0.44
1:B:647:CYS:SG	1:B:655:MET:HE3	2.58	0.43
1:A:587:LYS:HG3	3:A:108:HOH:O	2.17	0.43
1:B:617:MET:HB3	1:B:620:VAL:HB	1.99	0.43
1:A:628:ASN:O	1:A:629:ASN:C	2.56	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:421:ARG:HD2	1:B:444:PRO:HG3	2.00	0.43
1:A:419:ARG:NH1	1:A:493:ASN:O	2.52	0.43
1:A:419:ARG:HH12	1:A:494:PRO:N	2.16	0.43
1:B:501:LEU:C	1:B:501:LEU:HD13	2.38	0.43
1:B:493:ASN:ND2	1:B:493:ASN:N	2.63	0.43
1:B:417:ILE:HG21	1:B:441:TYR:CZ	2.53	0.43
1:A:419:ARG:HE	1:A:419:ARG:HB3	1.74	0.43
1:B:508:ARG:NE	3:B:115:HOH:O	2.46	0.43
1:B:485:LYS:NZ	1:B:487:ILE:HD13	2.34	0.42
1:B:426:ARG:NE	1:B:438:GLN:NE2	2.68	0.42
1:B:627:LYS:HB2	1:B:630:ASP:OD2	2.19	0.42
1:B:616:LEU:HD12	1:B:655:MET:HE1	2.02	0.42
1:A:459:CYS:HB2	3:A:74:HOH:O	2.19	0.42
1:B:501:LEU:HD12	1:B:503:THR:HG22	2.02	0.42
1:A:644:PRO:HB2	1:A:647:CYS:HB2	2.00	0.42
1:A:627:LYS:O	1:A:628:ASN:C	2.58	0.42
1:A:421:ARG:NH1	1:A:444:PRO:HD3	2.33	0.42
1:A:482:HIS:ND1	1:A:529:GLN:HB3	2.35	0.42
1:A:465:ARG:NH2	3:A:4:HOH:O	2.52	0.42
1:A:550:ARG:HD3	1:A:569:ARG:CZ	2.50	0.42
1:B:597:ARG:HG2	3:B:64:HOH:O	2.20	0.42
1:A:552:VAL:HG11	1:A:560:VAL:CG1	2.49	0.41
1:B:423:GLU:O	1:B:439:GLY:HA3	2.20	0.41
2:B:132:7PY:HAX3	2:B:132:7PY:CAN	2.50	0.41
1:A:444:PRO:C	1:A:446:ASN:N	2.72	0.41
1:A:500:GLU:O	2:A:131:7PY:HAB	2.20	0.41
1:B:508:ARG:HH11	1:B:508:ARG:CG	2.34	0.41
1:A:643:MET:HA	1:A:644:PRO:HD3	1.93	0.41
1:B:415:TYR:HE1	1:B:476:ARG:HG3	1.84	0.41
1:A:507:LEU:O	1:A:511:LEU:HG	2.20	0.41
1:A:544:HIS:O	1:A:545:ARG:CB	2.69	0.41
1:A:422:ILE:CD1	1:A:490:ILE:HD11	2.48	0.41
1:B:634:ARG:HG3	1:B:634:ARG:NH1	2.36	0.41
1:B:458:ASN:O	1:B:460:THR:N	2.51	0.41
1:B:438:GLN:O	1:B:438:GLN:HG3	2.21	0.41
1:A:583:LYS:HE3	1:A:583:LYS:HB2	1.96	0.41
1:B:602:ALA:HB3	3:B:117:HOH:O	2.20	0.41
1:B:424:LEU:N	1:B:424:LEU:HD23	2.35	0.41
2:A:131:7PY:HAN	2:A:131:7PY:HAX3	2.03	0.41
1:A:443:SER:OG	1:A:446:ASN:HB3	2.22	0.40
1:B:421:ARG:NH1	1:B:444:PRO:HG3	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:606:TRP:CE3	1:A:659:TRP:HA	2.57	0.40
1:B:423:GLU:O	1:B:439:GLY:CA	2.69	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	258/281 (92%)	234 (91%)	17 (7%)	7 (3%)	6	4
1	B	259/281 (92%)	240 (93%)	14 (5%)	5 (2%)	10	8
All	All	517/562 (92%)	474 (92%)	31 (6%)	12 (2%)	8	6

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	444	PRO
1	A	445	GLU
1	B	414	ASP
1	B	439	GLY
1	A	629	ASN
1	B	426	ARG
1	B	459	CYS
1	A	500	GLU
1	A	492	GLU
1	A	545	ARG
1	A	628	ASN
1	B	425	GLY

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	233 /248 (94%)	219 (94%)	14 (6%)	24	31
1	B	233 /248 (94%)	216 (93%)	17 (7%)	17	22
All	All	466 /496 (94%)	435 (93%)	31 (7%)	20	26

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	419	ARG
1	A	420	GLU
1	A	424	LEU
1	A	440	ILE
1	A	444	PRO
1	A	449	LEU
1	A	555	SER
1	A	597	ARG
1	A	598	ARG
1	A	624	GLN
1	A	637	ASN
1	A	651	LEU
1	A	653	SER
1	A	681	GLU
1	B	415	TYR
1	B	417	ILE
1	B	420	GLU
1	B	426	ARG
1	B	440	ILE
1	B	449	LEU
1	B	460	THR
1	B	469	LEU
1	B	470	GLN
1	B	493	ASN
1	B	508	ARG
1	B	515	LYS
1	B	517	SER

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Mol	Chain	Res	Type
1	B	555	SER
1	B	562	LEU
1	B	651	LEU
1	B	680	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	418	GLN
1	A	437	HIS
1	A	458	ASN
1	B	418	GLN
1	B	438	GLN
1	B	470	GLN
1	B	493	ASN
1	B	629	ASN
1	B	675	GLN

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 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	7PY	A	131	-	30,31,31	2.21	8 (26%)	33,43,43	2.59	13 (39%)
2	7PY	B	132	-	30,31,31	2.21	8 (26%)	33,43,43	2.54	13 (39%)

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	7PY	A	131	-	-	0/14/14/14	0/4/4/4
2	7PY	B	132	-	-	0/14/14/14	0/4/4/4

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	131	7PY	CAK-NAI	-4.89	1.33	1.39
2	B	132	7PY	CAK-NAI	-4.86	1.33	1.39
2	A	131	7PY	CAJ-NAI	-3.52	1.33	1.44
2	B	132	7PY	CAJ-NAI	-3.51	1.33	1.44
2	B	132	7PY	CAF-NAE	-3.27	1.33	1.40
2	A	131	7PY	CAF-NAE	-3.25	1.33	1.40
2	B	132	7PY	CAP-NAQ	2.78	1.40	1.34
2	A	131	7PY	CAP-NAQ	2.79	1.40	1.34
2	B	132	7PY	CAH-NAG	3.05	1.40	1.35
2	A	131	7PY	CAH-NAG	3.06	1.40	1.35
2	B	132	7PY	CAD-NAC	4.56	1.40	1.34
2	A	131	7PY	CAD-NAC	4.62	1.40	1.34
2	B	132	7PY	CAJ-NAQ	4.67	1.40	1.33
2	A	131	7PY	CAJ-NAQ	4.69	1.40	1.33
2	A	131	7PY	CAB-NAC	5.02	1.40	1.32
2	B	132	7PY	CAB-NAC	5.05	1.40	1.32

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	131	7PY	NAC-CAD-NAG	-7.00	118.88	126.62
2	B	132	7PY	NAC-CAD-NAG	-6.87	119.02	126.62
2	B	132	7PY	CBB-OBA-CAU	-4.13	111.28	117.54
2	A	131	7PY	CBB-OBA-CAU	-3.86	111.68	117.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	131	7PY	CAK-NAI-CAJ	-3.61	118.56	125.16
2	A	131	7PY	CAX-OAW-CAS	-3.38	112.42	117.54
2	B	132	7PY	CAX-OAW-CAS	-3.28	112.56	117.54
2	B	132	7PY	CAK-NAI-CAJ	-2.97	119.73	125.16
2	B	132	7PY	OAW-CAS-CAR	-2.95	119.17	124.21
2	B	132	7PY	OBA-CAU-CAV	-2.86	119.32	124.21
2	A	131	7PY	OBA-CAU-CAV	-2.80	119.43	124.21
2	A	131	7PY	OAW-CAS-CAR	-2.51	119.92	124.21
2	A	131	7PY	CAO-CAP-NAQ	-2.39	119.44	123.44
2	B	132	7PY	CAO-CAP-NAQ	-2.38	119.46	123.44
2	B	132	7PY	CAA-CAB-NAC	-2.37	119.72	124.06
2	A	131	7PY	CAA-CAB-NAC	-2.34	119.77	124.06
2	A	131	7PY	CAB-NAC-CAD	2.60	120.81	115.97
2	B	132	7PY	OAW-CAS-CAT	2.64	120.03	115.26
2	B	132	7PY	CAB-NAC-CAD	2.65	120.88	115.97
2	A	131	7PY	OAW-CAS-CAT	2.65	120.05	115.26
2	A	131	7PY	OBA-CAU-CAT	2.90	120.49	115.26
2	B	132	7PY	OBA-CAU-CAT	3.05	120.76	115.26
2	B	132	7PY	CAD-NAG-CAH	3.92	119.81	115.09
2	A	131	7PY	CAD-NAG-CAH	4.27	120.23	115.09
2	B	132	7PY	CAN-CAM-CAJ	7.05	121.41	117.06
2	A	131	7PY	CAN-CAM-CAJ	7.42	121.64	117.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	131	7PY	7	0
2	B	132	7PY	9	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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.