



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:00 PM GMT

PDB ID : 3VS0
Title : Crystal structure of HCK complexed with a pyrrolo-pyrimidine inhibitor N-[4-(4-amino-7-cyclopentyl-7H-pyrrolo[2,3-d]pyrimidin-5-yl)phenyl]benzamide
Authors : Kuratani, M.; Honda, K.; Tomabeche, Y.; Handa, N.; Yokoyama, S.
Deposited on : 2012-04-21
Resolution : 2.93 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

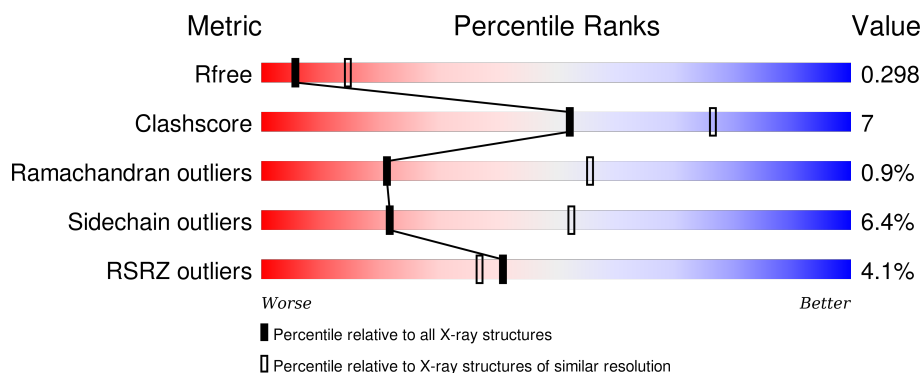
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.93 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	2057 (2.98-2.90)
Clashscore	102246	2308 (2.98-2.90)
Ramachandran outliers	100387	2245 (2.98-2.90)
Sidechain outliers	100360	2247 (2.98-2.90)
RSRZ outliers	91569	2065 (2.98-2.90)

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	454	<div> <div>7%</div> <div>78%</div> <div>15%</div> <div>6%</div> </div>
1	B	454	<div> <div>%</div> <div>73%</div> <div>21%</div> <div>.</div> </div>

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
3	CL	A	603	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7077 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 Tyrosine-protein kinase HCK.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	429	Total	C	N	O	P	S	0	0	0
			3463	2215	582	645	1	20			
1	B	434	Total	C	N	O	P	S	0	0	0
			3496	2233	588	654	1	20			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	78	GLY	-	EXPRESSION TAG	UNP P08631
A	79	ALA	-	EXPRESSION TAG	UNP P08631
A	80	MET	-	EXPRESSION TAG	UNP P08631
A	81	GLY	-	EXPRESSION TAG	UNP P08631
A	82	SER	-	EXPRESSION TAG	UNP P08631
A	83	GLY	-	EXPRESSION TAG	UNP P08631
A	84	ILE	-	EXPRESSION TAG	UNP P08631
A	85	ARG	-	EXPRESSION TAG	UNP P08631
A	528	GLU	GLN	ENGINEERED MUTATION	UNP P08631
A	529	GLU	GLN	ENGINEERED MUTATION	UNP P08631
A	530	ILE	GLN	ENGINEERED MUTATION	UNP P08631
B	78	GLY	-	EXPRESSION TAG	UNP P08631
B	79	ALA	-	EXPRESSION TAG	UNP P08631
B	80	MET	-	EXPRESSION TAG	UNP P08631
B	81	GLY	-	EXPRESSION TAG	UNP P08631
B	82	SER	-	EXPRESSION TAG	UNP P08631
B	83	GLY	-	EXPRESSION TAG	UNP P08631
B	84	ILE	-	EXPRESSION TAG	UNP P08631
B	85	ARG	-	EXPRESSION TAG	UNP P08631
B	528	GLU	GLN	ENGINEERED MUTATION	UNP P08631
B	529	GLU	GLN	ENGINEERED MUTATION	UNP P08631
B	530	ILE	GLN	ENGINEERED MUTATION	UNP P08631

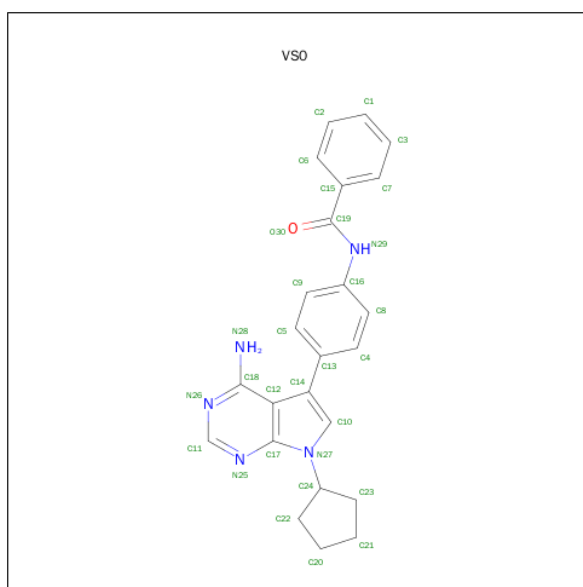
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ca 1 1	0	0
2	A	1	Total Ca 1 1	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0
3	A	1	Total Cl 1 1	0	0

- Molecule 4 is N-[4-(4-AMINO-7-CYCLOPENTYL-7H-PYRROLO[2,3-D]PYRIMIDIN-5-YL)PHENYL]BENZAMIDE (three-letter code: VS0) (formula: C₂₄H₂₃N₅O).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O 30 24 5 1	0	0
4	B	1	Total C N O 30 24 5 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	26	Total O 26 26	0	0

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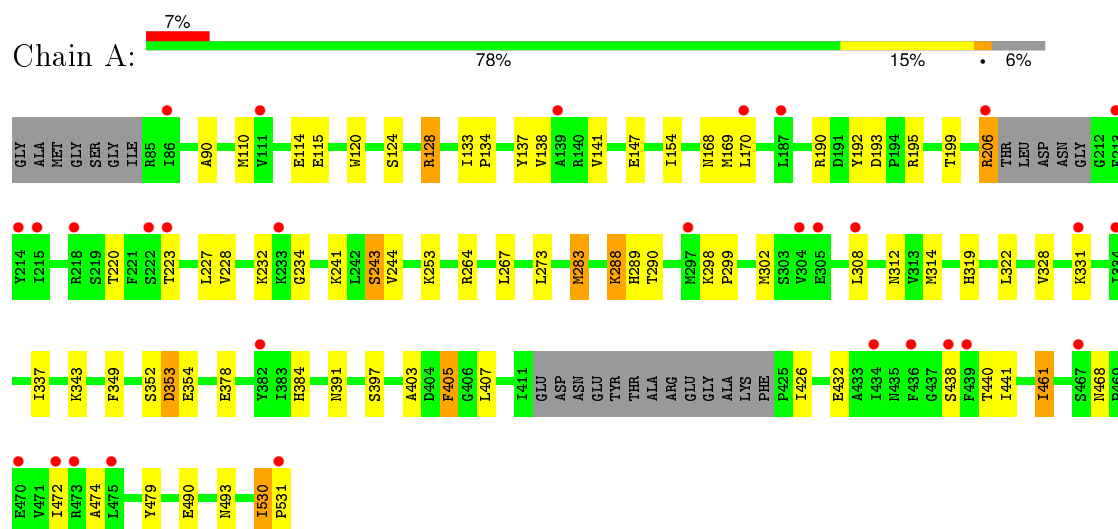
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	28	Total	O	0	0
			28	28		

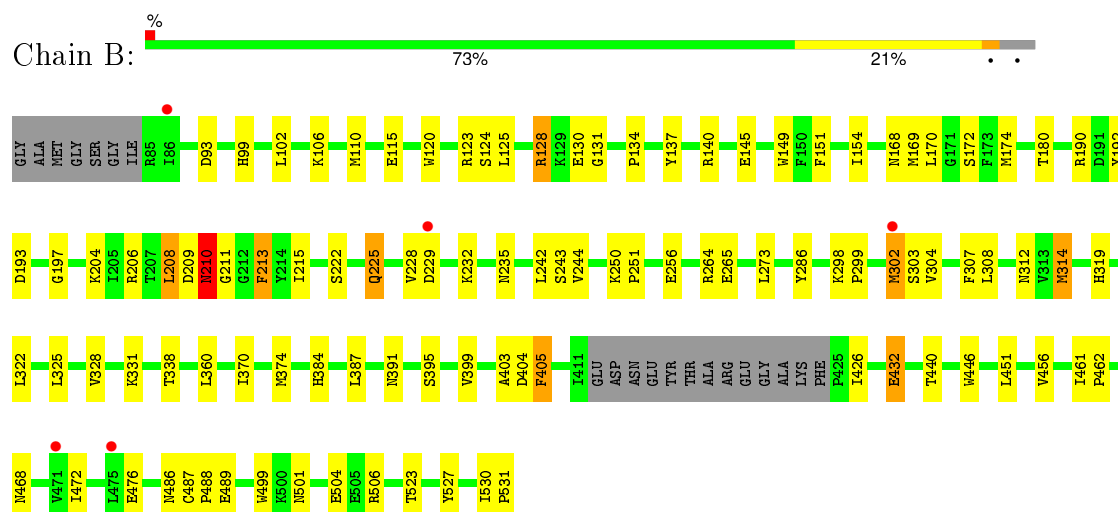
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Tyrosine-protein kinase HCK



- Molecule 1: Tyrosine-protein kinase HCK



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.72Å 96.05Å 181.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.29 – 2.93 39.29 – 2.93	Depositor EDS
% Data completeness (in resolution range)	99.6 (39.29-2.93) 99.6 (39.29-2.93)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 2.95Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.237 , 0.295 0.234 , 0.298	Depositor DCC
R_{free} test set	1411 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	72.7	Xtrriage
Anisotropy	0.137	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.7	EDS
Estimated twinning fraction	No twinning to report.	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Outliers	4 of 28235 reflections (0.014%)	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7077	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.22 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.1651e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: VS0, CA, PTR, CL

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	0.22	0/3528	0.40	0/4760
1	B	0.22	0/3562	0.42	1/4808 (0.0%)
All	All	0.22	0/7090	0.41	1/9568 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	210	ASN	N-CA-C	5.70	126.38	111.00

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	3463	0	3430	35	0
1	B	3496	0	3455	56	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	1	0
3	B	1	0	0	1	0
4	A	30	0	23	1	0
4	B	30	0	23	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	26	0	0	2	0
5	B	28	0	0	0	0
All	All	7077	0	6931	94	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 7.

All (94) 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:206:ARG:NH1	5:A:707:HOH:O	2.06	0.77
1:A:243:SER:OG	1:A:244:VAL:N	2.24	0.70
1:A:232:LYS:O	1:A:241:LYS:NZ	2.27	0.63
4:B:602:VS0:H9	4:B:602:VS0:H11	1.64	0.63
1:B:210:ASN:HB3	1:B:211:GLY:HA2	1.80	0.63
1:A:288:LYS:O	1:A:290:THR:N	2.30	0.62
4:B:602:VS0:C5	4:B:602:VS0:H11	2.13	0.62
1:A:128:ARG:NH2	3:A:603:CL:CL	2.71	0.60
1:B:125:LEU:HD23	1:B:128:ARG:HH21	1.67	0.60
1:B:264:ARG:NH2	1:B:331:LYS:O	2.34	0.58
1:B:456:VAL:HG11	1:B:487:CYS:HB2	1.86	0.56
1:B:215:ILE:HG22	1:B:530:ILE:HG21	1.87	0.56
1:B:180:THR:OG1	1:B:204:LYS:NZ	2.37	0.56
1:A:147:GLU:OE2	5:A:713:HOH:O	2.18	0.56
1:B:319:HIS:HB3	1:B:322:LEU:HG	1.90	0.54
1:B:264:ARG:HH21	1:B:331:LYS:HB2	1.71	0.54
1:B:530:ILE:HD12	1:B:531:PRO:HD2	1.89	0.54
1:A:319:HIS:HB3	1:A:322:LEU:HG	1.91	0.53
1:B:243:SER:OG	1:B:244:VAL:N	2.38	0.53
1:A:264:ARG:NH2	1:A:331:LYS:O	2.39	0.52
1:B:149:TRP:HZ2	1:B:225:GLN:HG2	1.73	0.52
1:B:426:ILE:HD11	1:B:468:ASN:HB3	1.90	0.52
1:A:343:LYS:NZ	1:A:354:GLU:OE2	2.43	0.52
1:B:395:SER:HB3	1:B:399:VAL:H	1.76	0.51
1:A:391:ASN:HB3	1:A:403:ALA:O	2.11	0.50
1:B:232:LYS:HG2	1:B:242:LEU:HB2	1.92	0.50
1:A:267:LEU:HD11	1:A:337:ILE:HD11	1.94	0.50
1:A:273:LEU:HD11	1:A:283:MET:HB3	1.93	0.49
1:B:209:ASP:HA	1:B:210:ASN:C	2.32	0.49
1:B:204:LYS:NZ	1:B:527:PTR:O3P	2.40	0.49
1:B:265:GLU:N	1:B:265:GLU:OE1	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:PHE:HB3	1:B:154:ILE:HD13	1.94	0.49
1:A:168:ASN:HB2	1:A:190:ARG:HH11	1.78	0.48
1:B:228:VAL:O	1:B:232:LYS:HG3	2.14	0.48
1:B:325:LEU:HA	1:B:338:THR:HA	1.96	0.48
1:A:170:LEU:HD11	1:A:192:TYR:HB2	1.96	0.48
1:B:210:ASN:CB	1:B:211:GLY:HA2	2.43	0.48
1:B:93:ASP:OD1	1:B:106:LYS:N	2.44	0.48
1:A:426:ILE:HD11	1:A:468:ASN:HB3	1.94	0.48
1:B:314:MET:N	1:B:314:MET:SD	2.86	0.48
1:B:273:LEU:HB3	4:B:602:VS0:H16	1.95	0.47
1:B:128:ARG:NH2	3:B:603:CL:CL	2.68	0.47
1:B:102:LEU:HG	1:B:131:GLY:HA3	1.96	0.47
1:A:90:ALA:HA	1:A:138:VAL:HG12	1.97	0.47
1:B:208:LEU:O	1:B:211:GLY:HA3	2.15	0.47
1:B:489:GLU:H	1:B:489:GLU:CD	2.17	0.47
1:A:115:GLU:HA	1:A:120:TRP:CD1	2.51	0.46
1:A:352:SER:OG	1:A:353:ASP:N	2.49	0.46
1:A:193:ASP:OD1	1:A:195:ARG:NE	2.39	0.46
1:A:378:GLU:HG3	1:A:441:ILE:HG12	1.97	0.45
1:A:110:MET:SD	1:A:133:ILE:HD13	2.57	0.45
1:B:374:MET:HG3	1:B:387:LEU:HD11	1.98	0.45
1:B:110:MET:HB3	1:B:124:SER:HA	1.99	0.45
1:B:170:LEU:HD21	1:B:192:TYR:HB2	1.99	0.45
1:A:384:HIS:HD2	1:A:405:PHE:HB2	1.81	0.45
4:A:602:VS0:H11	4:A:602:VS0:C5	2.30	0.44
1:B:432:GLU:OE2	1:B:506:ARG:NH1	2.49	0.44
1:A:228:VAL:O	1:A:232:LYS:HG3	2.18	0.44
1:A:110:MET:HB3	1:A:124:SER:HA	1.99	0.44
1:A:530:ILE:HA	1:A:531:PRO:HD3	1.70	0.43
1:A:234:GLY:H	1:A:241:LYS:HD2	1.84	0.43
1:B:486:ASN:N	1:B:486:ASN:OD1	2.48	0.43
1:B:250:LYS:HA	1:B:251:PRO:HD3	1.91	0.43
4:B:602:VS0:H9	4:B:602:VS0:N28	2.29	0.43
1:A:474:ALA:HB1	1:A:479:TYR:HB3	2.00	0.43
1:B:168:ASN:HB2	1:B:190:ARG:HH11	1.83	0.43
1:B:206:ARG:O	1:B:213:PHE:HA	2.19	0.43
1:B:256:GLU:OE2	1:B:286:TYR:OH	2.30	0.43
1:A:298:LYS:HA	1:A:299:PRO:HD3	1.89	0.43
1:B:123:ARG:HB2	1:B:130:GLU:HG2	2.01	0.42
1:A:349:PHE:O	1:A:352:SER:HB3	2.18	0.42
1:B:432:GLU:HG3	1:B:432:GLU:H	1.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:432:GLU:O	1:A:438:SER:OG	2.33	0.42
1:A:134:PRO:HG2	1:A:137:TYR:HB2	2.01	0.42
1:B:151:PHE:CD1	1:B:174:MET:HB2	2.54	0.42
1:A:468:ASN:O	1:A:472:ILE:HG13	2.20	0.42
1:A:461:ILE:H	1:A:461:ILE:HG13	1.55	0.41
1:B:225:GLN:NE2	1:B:229:ASP:OD1	2.52	0.41
1:B:370:ILE:HD13	1:B:451:LEU:HD21	2.01	0.41
1:B:134:PRO:HG2	1:B:137:TYR:HB2	2.02	0.41
1:B:115:GLU:HA	1:B:120:TRP:CD1	2.55	0.41
1:A:384:HIS:CD2	1:A:405:PHE:HB2	2.55	0.41
1:B:123:ARG:NH1	1:B:130:GLU:OE2	2.54	0.41
1:B:461:ILE:HA	1:B:462:PRO:HD3	1.92	0.41
1:B:172:SER:HA	1:B:244:VAL:O	2.20	0.41
1:B:468:ASN:O	1:B:472:ILE:HG13	2.20	0.41
1:B:384:HIS:CD2	1:B:405:PHE:HB2	2.56	0.41
1:B:391:ASN:HB3	1:B:403:ALA:O	2.21	0.41
1:B:235:ASN:N	1:B:235:ASN:OD1	2.53	0.41
1:A:490:GLU:HG2	1:B:523:THR:O	2.21	0.41
1:B:446:TRP:CE3	1:B:499:TRP:HA	2.57	0.40
1:B:168:ASN:O	1:B:190:ARG:NH1	2.55	0.40
1:B:360:LEU:HD21	1:B:488:PRO:HG3	2.03	0.40
1:B:298:LYS:HA	1:B:299:PRO:HD3	1.85	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	422/454 (93%)	402 (95%)	17 (4%)	3 (1%)	26	62
1	B	429/454 (94%)	401 (94%)	23 (5%)	5 (1%)	16	47
All	All	851/908 (94%)	803 (94%)	40 (5%)	8 (1%)	21	56

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	289	HIS
1	B	210	ASN
1	B	197	GLY
1	A	288	LYS
1	A	405	PHE
1	B	302	MET
1	B	405	PHE
1	B	304	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	374/393 (95%)	349 (93%)	25 (7%)	20	50
1	B	377/393 (96%)	354 (94%)	23 (6%)	23	55
All	All	751/786 (96%)	703 (94%)	48 (6%)	22	52

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	GLU
1	A	128	ARG
1	A	141	VAL
1	A	154	ILE
1	A	169	MET
1	A	199	THR
1	A	206	ARG
1	A	220	THR
1	A	223	THR
1	A	227	LEU
1	A	243	SER
1	A	253	LYS
1	A	283	MET
1	A	302	MET
1	A	308	LEU

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Mol	Chain	Res	Type
1	A	312	ASN
1	A	314	MET
1	A	328	VAL
1	A	353	ASP
1	A	397	SER
1	A	407	LEU
1	A	440	THR
1	A	461	ILE
1	A	493	ASN
1	A	530	ILE
1	B	99	HIS
1	B	128	ARG
1	B	140	ARG
1	B	145	GLU
1	B	169	MET
1	B	193	ASP
1	B	208	LEU
1	B	213	PHE
1	B	222	SER
1	B	225	GLN
1	B	302	MET
1	B	303	SER
1	B	307	PHE
1	B	308	LEU
1	B	312	ASN
1	B	314	MET
1	B	328	VAL
1	B	404	ASP
1	B	432	GLU
1	B	440	THR
1	B	476	GLU
1	B	501	ASN
1	B	504	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	277	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	PTR	A	527	1,2	14,16,17	1.24	1 (7%)	18,22,24	0.61	0
1	PTR	B	527	1	14,16,17	1.25	1 (7%)	18,22,24	0.64	0

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
1	PTR	A	527	1,2	-	0/9/11/13	0/1/1/1
1	PTR	B	527	1	-	0/9/11/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	527	PTR	OH-CZ	-4.08	1.30	1.40
1	A	527	PTR	OH-CZ	-4.06	1.30	1.40

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	527	PTR	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
4	VS0	A	602	-	32,34,34	0.72	1 (3%)	35,48,48	1.90	2 (5%)
4	VS0	B	602	-	32,34,34	0.67	1 (3%)	35,48,48	2.01	2 (5%)

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
4	VS0	A	602	-	-	0/12/23/23	0/5/5/5
4	VS0	B	602	-	-	0/12/23/23	0/5/5/5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	602	VS0	C10-N27	-2.77	1.34	1.38
4	B	602	VS0	C10-N27	-2.55	1.34	1.38

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	602	VS0	N25-C11-N26	-10.47	120.88	128.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	602	VS0	N25-C11-N26	-9.93	121.29	128.89
4	B	602	VS0	C22-C24-C23	2.75	107.16	104.32
4	A	602	VS0	C22-C24-C23	2.85	107.27	104.32

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	602	VS0	1	0
4	B	602	VS0	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 ⓘ

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, 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	428/454 (94%)	0.53	30 (7%) 19 15	47, 80, 110, 135	0
1	B	433/454 (95%)	0.33	5 (1%) 81 80	49, 76, 101, 124	1 (0%)
All	All	861/908 (94%)	0.43	35 (4%) 41 37	47, 78, 107, 135	1 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	475	LEU	4.3
1	A	434	ILE	3.5
1	A	305	GLU	3.3
1	A	308	LEU	3.1
1	A	213	PHE	3.1
1	A	222	SER	3.1
1	A	214	TYR	3.0
1	A	334	ILE	2.8
1	A	531	PRO	2.8
1	A	86	ILE	2.7
1	B	475	LEU	2.6
1	A	473	ARG	2.6
1	B	86	ILE	2.5
1	A	233	LYS	2.4
1	A	472	ILE	2.4
1	A	438	SER	2.4
1	A	436	PHE	2.4
1	A	382	TYR	2.4
1	A	304	VAL	2.4
1	A	111	VAL	2.2
1	A	439	PHE	2.2
1	A	297	MET	2.2
1	A	215	ILE	2.2
1	A	470	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	218	ARG	2.2
1	A	467	SER	2.2
1	A	331	LYS	2.2
1	A	187	LEU	2.1
1	B	302	MET	2.1
1	A	206	ARG	2.1
1	A	223	THR	2.1
1	A	139	ALA	2.1
1	B	471	VAL	2.1
1	B	229	ASP	2.0
1	A	170	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	PTR	B	527	16/17	0.93	0.20	-	54,63,74,75	0
1	PTR	A	527	16/17	0.93	0.19	-	55,67,77,77	0

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(Å ²)	Q<0.9
3	CL	A	603	1/1	0.87	0.46	3.52	92,92,92,92	0
4	VS0	A	602	30/30	0.95	0.28	1.03	59,67,87,93	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	VS0	B	602	30/30	0.96	0.26	0.48	50,60,81,85	0
3	CL	B	603	1/1	0.91	0.29	0.45	84,84,84,84	0
2	CA	A	601	1/1	0.73	0.25	-	82,82,82,82	0
2	CA	B	601	1/1	0.77	0.31	-	70,70,70,70	0

6.5 Other polymers [i](#)

There are no such residues in this entry.