



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:00 PM GMT

PDB ID : 3VS2
Title : Crystal structure of HCK complexed with a pyrrolo-pyrimidine inhibitor 7-[cis-4-(4-methylpiperazin-1-yl)cyclohexyl]-5-(4-phenoxyphenyl)-7H-pyrrolo[2,3-d]pyrimidin-4-amine
Authors : Kuratani, M.; Tomabechei, Y.; Niwa, H.; Parker, J.L.; Handa, N.; Yokoyama, S.
Deposited on : 2012-04-21
Resolution : 2.61 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

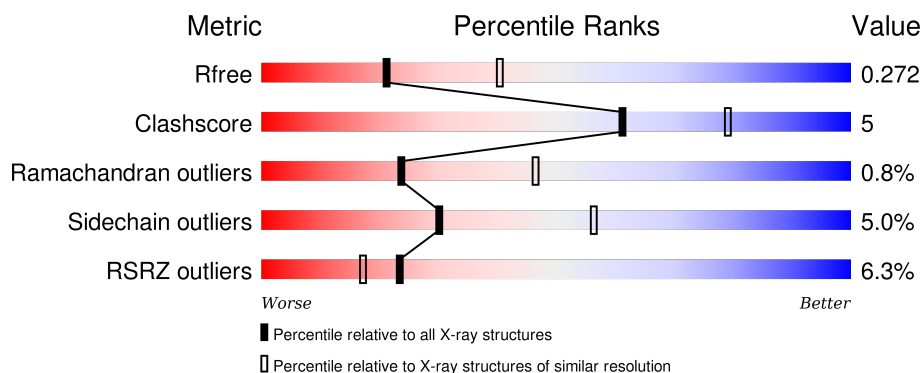
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	 6% 80% 13% • 5%
1	B	454	 6% 79% 16% • •

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7189 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	431	Total	C	N	O	P	S	0	0	0
			3487	2231	587	648	1	20			
1	B	434	Total	C	N	O	P	S	0	0	0
			3504	2238	591	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 7-[CIS-4-(4-METHYLPIPERAZIN-1-YL)CYCLOHEXYL]-5-(4-PHENOXYPHENYL)-7H-PYRROLO[2,3-D]PYRIMIDIN-4-AMINE (three-letter code: VSB) (formula:

The chemical structure of VSB (Vincristine Sulfate) is a complex polycyclic alkaloid. It features a central tetracyclic core with several fused and linked rings. Key components include a dihydroquinoline ring system, a tetrahydroisoquinoline ring, and a complex polycyclic system with multiple nitrogen atoms. The structure is highly substituted with various functional groups, including a sulfate group (SO₃Na) and several hydroxyl groups. The overall structure is shown in a 2D representation with various atoms labeled with their respective symbols (C, H, N, O, S, Na).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 36	C 29	N 6	O 1	0	0
2	B	1	Total 36	C 29	N 6	O 1	0	0

- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|-----------------|---------|---------|
| 3 | B | 1 | Total Ca
1 1 | 0 | 0 |
| 3 | A | 1 | Total Ca
1 1 | 0 | 0 |

- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|-----------------|---------|---------|
| 4 | B | 1 | Total Cl
1 1 | 0 | 0 |
| 4 | A | 1 | Total Cl
1 1 | 0 | 0 |

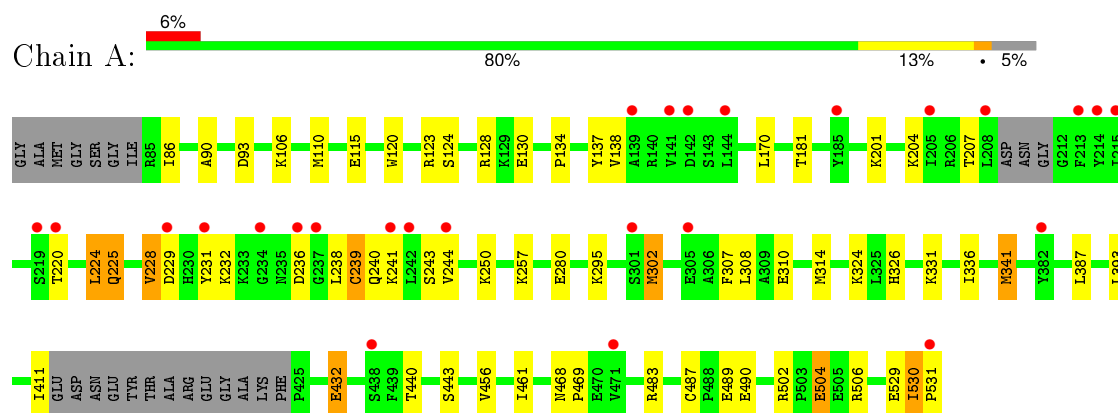
- 

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	67	Total 67	O 67	0	0
5	B	55	Total 55	O 55	0	0

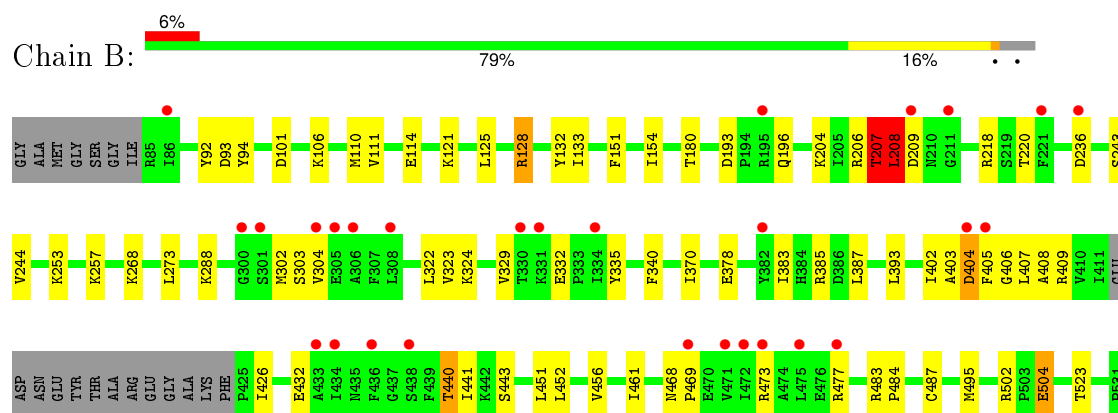
3 Residue-property plots

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: Tyrosine-protein kinase HCK



• Molecule 1: Tyrosine-protein kinase HCK



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.51Å 73.09Å 180.96Å 90.00° 96.26° 90.00°	Depositor
Resolution (Å)	40.69 – 2.61 40.69 – 2.61	Depositor EDS
% Data completeness (in resolution range)	98.4 (40.69-2.61) 98.5 (40.69-2.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.226 , 0.279 0.221 , 0.272	Depositor DCC
R_{free} test set	1903 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	62.8	Xtriage
Anisotropy	0.585	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 59.1	EDS
Estimated twinning fraction	0.039 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	3 of 38045 reflections (0.008%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7189	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: VSB, CL, CA, PTR

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	0/3552	0.40	0/4792
1	B	0.22	0/3570	0.40	0/4818
All	All	0.22	0/7122	0.40	0/9610

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	3487	0	3468	33	0
1	B	3504	0	3473	41	0
2	A	36	0	31	3	0
2	B	36	0	32	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	1	0
5	A	67	0	0	1	0
5	B	55	0	0	3	0
All	All	7189	0	7004	77	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 5.

All (77) 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:B:128:ARG:NH2	4:B:603:CL:CL	2.42	0.88
1:B:180:THR:OG1	1:B:204:LYS:NZ	2.30	0.65
1:B:323:VAL:HB	1:B:403:ALA:HB2	1.80	0.64
1:B:483:ARG:NH1	1:B:487:CYS:O	2.30	0.63
1:B:406:GLY:O	1:B:408:ALA:N	2.33	0.61
1:B:114:GLU:HB3	1:B:121:LYS:HB3	1.81	0.61
1:B:404:ASP:HA	1:B:405:PHE:HB3	1.82	0.60
1:A:225:GLN:NE2	1:A:229:ASP:OD2	2.33	0.60
1:A:530:ILE:HG12	1:A:531:PRO:HD2	1.82	0.60
1:A:502:ARG:HB3	1:A:504:GLU:HG2	1.83	0.59
1:B:385:ARG:NH1	1:B:408:ALA:O	2.35	0.59
2:B:601:VSF:H24	2:B:601:VSF:CAL	2.15	0.59
2:A:601:VSF:CAL	2:A:601:VSF:H24	2.17	0.56
1:A:302:MET:N	1:A:302:MET:SD	2.75	0.56
1:A:432:GLU:OE2	1:A:506:ARG:NH2	2.28	0.56
1:B:502:ARG:HG2	1:B:504:GLU:HG2	1.87	0.55
1:A:93:ASP:O	1:A:250:LYS:NZ	2.40	0.55
1:B:236:ASP:OD2	5:B:747:HOH:O	2.17	0.54
1:B:151:PHE:HB3	1:B:154:ILE:HG13	1.90	0.53
1:B:383:ILE:HD11	1:B:409:ARG:HD2	1.92	0.51
1:B:110:MET:HA	1:B:125:LEU:HG	1.92	0.51
1:A:181:THR:OG1	1:A:204:LYS:NZ	2.40	0.50
1:A:483:ARG:NH1	1:A:487:CYS:O	2.43	0.48
2:B:601:VSF:H24	2:B:601:VSF:H28	1.78	0.48
1:A:243:SER:OG	1:A:244:VAL:N	2.47	0.48
1:B:268:LYS:HE3	1:B:288:LYS:HE3	1.96	0.48
1:B:207:THR:O	1:B:208:LEU:HB2	2.14	0.47
1:A:490:GLU:HG2	1:B:523:THR:O	2.15	0.47
1:B:403:ALA:HA	1:B:405:PHE:HB3	1.96	0.47
1:B:111:VAL:HG23	1:B:125:LEU:HD21	1.97	0.47
1:A:224:LEU:O	1:A:228:VAL:HG12	2.15	0.47
1:B:370:ILE:HD13	1:B:451:LEU:HD21	1.97	0.46
1:B:323:VAL:HG21	1:B:393:LEU:HD12	1.96	0.46
1:B:253:LYS:HB3	1:B:253:LYS:HE2	1.70	0.46
1:B:473:ARG:O	1:B:477:ARG:HG3	2.15	0.46
1:B:461:ILE:HD12	1:B:461:ILE:H	1.80	0.46
1:B:93:ASP:OD1	1:B:106:LYS:N	2.42	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:452:LEU:HD23	1:B:495:MET:HG3	1.98	0.46
1:A:231:TYR:HE1	1:A:236:ASP:HB2	1.81	0.46
1:B:378:GLU:HG3	1:B:441:ILE:HG12	1.97	0.46
1:A:90:ALA:HA	1:A:138:VAL:HG12	1.98	0.45
1:B:236:ASP:O	5:B:747:HOH:O	2.21	0.45
1:A:123:ARG:NH1	1:A:130:GLU:OE2	2.50	0.45
1:A:123:ARG:HB2	1:A:130:GLU:HG2	1.98	0.45
1:A:93:ASP:OD1	1:A:106:LYS:N	2.42	0.44
1:B:322:LEU:HD22	1:B:402:ILE:HB	1.98	0.44
1:A:110:MET:HG3	1:A:124:SER:HA	1.99	0.44
1:A:201:LYS:HG3	1:A:239:CYS:SG	2.58	0.44
1:B:273:LEU:HD11	1:B:340:PHE:HE1	1.83	0.44
1:B:426:ILE:HD11	1:B:468:ASN:HB3	1.99	0.44
1:B:110:MET:SD	1:B:133:ILE:HD13	2.58	0.43
1:A:489:GLU:CD	1:B:206:ARG:HH12	2.22	0.43
1:A:232:LYS:HB3	1:A:232:LYS:HE2	1.84	0.43
1:B:94:TYR:HE1	1:B:253:LYS:HD2	1.83	0.43
1:B:393:LEU:HG	5:B:731:HOH:O	2.19	0.42
1:A:170:LEU:HD12	1:A:240:GLN:HG3	2.01	0.42
1:B:440:THR:O	1:B:443:SER:OG	2.21	0.42
2:B:601:VSB:H10	2:B:601:VSB:H36	1.71	0.42
1:A:307:PHE:CZ	1:A:336:ILE:HD11	2.55	0.42
1:B:101:ASP:HA	1:B:132:TYR:H	1.85	0.41
1:B:243:SER:OG	1:B:244:VAL:N	2.52	0.41
1:A:341:MET:O	2:A:601:VSB:H22	2.20	0.41
2:A:601:VSB:H24	2:A:601:VSB:H28	1.82	0.41
1:A:341:MET:HG2	1:A:393:LEU:HB3	2.02	0.41
1:B:468:ASN:HB2	1:B:469:PRO:HD3	2.02	0.41
1:A:295:LYS:NZ	5:A:712:HOH:O	2.54	0.41
1:A:461:ILE:HD12	1:A:461:ILE:H	1.85	0.41
1:A:115:GLU:HA	1:A:120:TRP:CD1	2.56	0.41
1:B:324:LYS:HE3	1:B:324:LYS:HB3	1.86	0.41
1:B:483:ARG:HA	1:B:484:PRO:HD2	1.92	0.41
1:A:324:LYS:HD2	1:A:326:HIS:CE1	2.55	0.40
1:B:329:VAL:HB	1:B:335:TYR:HB2	2.03	0.40
1:A:468:ASN:HB2	1:A:469:PRO:HD3	2.03	0.40
1:A:310:GLU:O	1:A:314:MET:HG3	2.21	0.40
1:A:134:PRO:HG2	1:A:137:TYR:HB2	2.03	0.40
1:A:440:THR:O	1:A:443:SER:OG	2.37	0.40
1:A:411:ILE:H	1:A:411:ILE:HG13	1.67	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	424/454 (93%)	398 (94%)	24 (6%)	2 (0%)	34	60
1	B	429/454 (94%)	398 (93%)	26 (6%)	5 (1%)	16	33
All	All	853/908 (94%)	796 (93%)	50 (6%)	7 (1%)	24	46

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	207	THR
1	B	208	LEU
1	B	407	LEU
1	A	239	CYS
1	B	303	SER
1	B	196	GLN
1	A	207	THR

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	378/393 (96%)	358 (95%)	20 (5%)	28	53
1	B	379/393 (96%)	361 (95%)	18 (5%)	32	59
All	All	757/786 (96%)	719 (95%)	38 (5%)	30	56

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86	ILE
1	A	128	ARG
1	A	220	THR
1	A	224	LEU
1	A	225	GLN
1	A	228	VAL
1	A	238	LEU
1	A	241	LYS
1	A	257	LYS
1	A	280	GLU
1	A	302	MET
1	A	308	LEU
1	A	331	LYS
1	A	341	MET
1	A	387	LEU
1	A	432	GLU
1	A	456	VAL
1	A	504	GLU
1	A	529	GLU
1	A	530	ILE
1	B	92	TYR
1	B	128	ARG
1	B	193	ASP
1	B	207	THR
1	B	208	LEU
1	B	209	ASP
1	B	218	ARG
1	B	220	THR
1	B	257	LYS
1	B	302	MET
1	B	304	VAL
1	B	332	GLU
1	B	387	LEU
1	B	404	ASP
1	B	432	GLU
1	B	440	THR
1	B	456	VAL
1	B	504	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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	1,3	14,16,17	1.31	1 (7%)	18,22,24	0.95	1 (5%)
1	PTR	B	527	1,3	14,16,17	1.17	1 (7%)	18,22,24	0.76	1 (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
1	PTR	A	527	1,3	-	0/9/11/13	0/1/1/1
1	PTR	B	527	1,3	-	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	A	527	PTR	OH-CZ	-4.34	1.30	1.40
1	B	527	PTR	OH-CZ	-4.19	1.30	1.40

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	527	PTR	O-C-CA	-3.29	116.91	125.49
1	B	527	PTR	O-C-CA	-2.50	118.97	125.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	$\# Z > 2$
2	VSB	A	601	-	39,41,41	3.01	12 (30%)	47,58,58	9.85	25 (53%)
2	VSB	B	601	-	39,41,41	2.71	14 (35%)	47,58,58	9.13	22 (46%)

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	VSB	A	601	-	-	0/12/36/36	1/6/6/6
2	VSB	B	601	-	-	0/12/36/36	1/6/6/6

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	VSB	CAR-CAS	-7.93	1.32	1.52
2	A	601	VSB	CAU-CAV	-7.17	1.34	1.52
2	B	601	VSB	CAU-CAV	-6.78	1.35	1.52
2	B	601	VSB	CBB-NAW	-5.75	1.35	1.47
2	A	601	VSB	CAX-NAW	-5.75	1.35	1.47

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	VS	CAR-CAS	-5.66	1.38	1.52
2	A	601	VS	CAH-NAG	-4.14	1.32	1.38
2	A	601	VS	CAI-CAJ	-4.04	1.42	1.49
2	B	601	VS	CAI-CAJ	-4.00	1.42	1.49
2	B	601	VS	CAH-NAG	-3.80	1.32	1.38
2	A	601	VS	CAH-CAI	-3.79	1.34	1.38
2	B	601	VS	CAH-CAI	-3.28	1.34	1.38
2	B	601	VS	CAY-NAZ	-2.14	1.41	1.46
2	B	601	VS	CAY-CAX	3.12	1.64	1.51
2	B	601	VS	CAS-CAT	3.16	1.61	1.52
2	B	601	VS	CAU-CAT	3.18	1.61	1.52
2	A	601	VS	CBA-CBB	3.58	1.66	1.51
2	B	601	VS	CAR-CAQ	3.82	1.62	1.52
2	B	601	VS	CAV-CAQ	3.83	1.62	1.52
2	A	601	VS	CAV-CAQ	3.94	1.62	1.52
2	A	601	VS	CBB-NAW	4.53	1.56	1.47
2	B	601	VS	CAX-NAW	4.75	1.57	1.47
2	A	601	VS	CAR-CAQ	5.35	1.66	1.52
2	A	601	VS	CAS-CAT	5.42	1.67	1.52
2	A	601	VS	CAU-CAT	5.50	1.67	1.52
2	B	601	VS	CBA-NAZ	5.85	1.59	1.46

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	VS	CAX-CAY-NAZ	-33.45	75.46	110.79
2	A	601	VS	CBB-CBA-NAZ	-30.97	78.07	110.79
2	A	601	VS	CAS-CAR-CAQ	-21.97	83.46	109.75
2	B	601	VS	CAR-CAQ-NAG	-20.56	75.15	110.76
2	A	601	VS	CAX-NAW-CBB	-20.24	72.54	109.23
2	B	601	VS	CAU-CAT-NAW	-20.21	67.06	112.57
2	B	601	VS	CAX-NAW-CBB	-18.81	75.14	109.23
2	A	601	VS	CAU-CAT-NAW	-16.72	74.91	112.57
2	A	601	VS	CAR-CAQ-NAG	-16.25	82.61	110.76
2	A	601	VS	CAX-CAY-NAZ	-15.57	94.34	110.79
2	A	601	VS	CAV-CAU-CAT	-14.81	78.82	109.54
2	B	601	VS	CAY-NAZ-CBA	-12.40	93.27	109.53
2	B	601	VS	CAS-CAR-CAQ	-8.51	99.57	109.75
2	B	601	VS	N3-C2-N1	-8.49	122.39	128.89
2	A	601	VS	N3-C2-N1	-7.57	123.10	128.89
2	B	601	VS	CAV-CAU-CAT	-6.89	95.25	109.54
2	B	601	VS	CBA-CBB-NAW	-4.85	99.43	110.79

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	VS	CAI-C5-C4	-4.73	103.76	107.54
2	A	601	VS	CAR-CAS-CAT	-4.56	100.09	109.54
2	B	601	VS	CAI-C5-C4	-3.53	104.72	107.54
2	A	601	VS	CAU-CAV-CAQ	-2.27	107.03	109.75
2	B	601	VS	OBD-CAN-CAO	2.03	126.68	119.42
2	A	601	VS	OBD-CAN-CAO	2.07	126.83	119.42
2	B	601	VS	CAU-CAT-CAS	2.63	117.42	111.22
2	B	601	VS	CBE-OBD-CAN	2.76	125.69	118.81
2	A	601	VS	CBE-OBD-CAN	2.84	125.90	118.81
2	A	601	VS	CBC-NAZ-CBA	2.93	115.20	110.63
2	A	601	VS	CBB-NAW-CAT	3.04	118.33	112.37
2	B	601	VS	CBB-CBA-NAZ	3.47	114.45	110.79
2	B	601	VS	CAH-NAG-CAQ	5.62	130.17	125.44
2	A	601	VS	CBC-NAZ-CAY	6.65	120.99	110.63
2	B	601	VS	CAU-CAV-CAQ	7.21	118.37	109.75
2	A	601	VS	CAS-CAT-NAW	7.32	129.05	112.57
2	A	601	VS	CAV-CAQ-NAG	7.64	123.99	110.76
2	A	601	VS	CAV-CAQ-CAR	7.94	125.42	110.13
2	A	601	VS	CAU-CAT-CAS	8.30	130.78	111.22
2	A	601	VS	CAH-NAG-CAQ	9.06	133.06	125.44
2	A	601	VS	CAY-NAZ-CBA	10.16	122.86	109.53
2	B	601	VS	CAS-CAT-NAW	10.51	136.25	112.57
2	B	601	VS	CAV-CAQ-NAG	12.08	131.69	110.76
2	B	601	VS	CAY-CAX-NAW	12.90	141.04	110.79
2	B	601	VS	CBC-NAZ-CBA	13.91	132.32	110.63
2	A	601	VS	CAY-CAX-NAW	14.77	145.42	110.79
2	B	601	VS	CBC-NAZ-CAY	15.04	134.08	110.63
2	A	601	VS	CBA-CBB-NAW	15.37	146.82	110.79
2	B	601	VS	CBB-NAW-CAT	16.13	143.98	112.37
2	A	601	VS	CAX-NAW-CAT	24.90	161.16	112.37

There are no chirality outliers.

There are no torsion outliers.

All (2) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	VS	CAX-CAY-CBA-CBB-NAW-NAZ
2	B	601	VS	CAX-CAY-CBA-CBB-NAW-NAZ

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	VS	3	0
2	B	601	VS	3	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	430/454 (94%)	0.53	26 (6%)	25 18	41, 70, 113, 134	0
1	B	433/454 (95%)	0.58	28 (6%)	22 16	46, 75, 114, 143	1 (0%)
All	All	863/908 (95%)	0.55	54 (6%)	23 17	41, 73, 114, 143	1 (0%)

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	VAL	9.9
1	B	475	LEU	6.1
1	B	471	VAL	5.2
1	A	236	ASP	5.1
1	B	211	GLY	4.7
1	A	139	ALA	4.2
1	A	531	PRO	4.2
1	A	241	LYS	4.0
1	B	305	GLU	3.8
1	A	229	ASP	3.6
1	A	215	ILE	3.5
1	B	301	SER	3.4
1	B	306	ALA	3.4
1	A	301	SER	3.4
1	B	434	ILE	3.4
1	B	438	SER	3.4
1	B	334	ILE	3.4
1	A	382	TYR	3.3
1	B	469	PRO	3.3
1	A	234	GLY	3.3
1	A	142	ASP	3.0
1	A	237	GLY	2.9
1	A	219	SER	2.9
1	A	213	PHE	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	242	LEU	2.7
1	A	205	ILE	2.6
1	B	209	ASP	2.6
1	B	382	TYR	2.6
1	A	471	VAL	2.5
1	A	231	TYR	2.5
1	B	473	ARG	2.5
1	A	144	LEU	2.4
1	B	330	THR	2.4
1	B	433	ALA	2.4
1	B	86	ILE	2.4
1	A	244	VAL	2.3
1	A	305	GLU	2.3
1	B	195	ARG	2.3
1	A	438	SER	2.3
1	A	214	TYR	2.3
1	A	141	VAL	2.3
1	A	220	THR	2.2
1	B	300	GLY	2.2
1	B	331	LYS	2.1
1	B	308	LEU	2.1
1	B	477	ARG	2.1
1	B	236	ASP	2.1
1	A	208	LEU	2.1
1	A	185	TYR	2.1
1	B	404	ASP	2.1
1	B	405	PHE	2.1
1	B	472	ILE	2.0
1	B	221	PHE	2.0
1	B	436	PHE	2.0

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

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	A	527	16/17	0.96	0.14	-	49,68,77,85	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	PTR	B	527	16/17	0.95	0.19	-	43,59,64,71	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(\AA^2)	Q<0.9
2	VSB	B	601	36/36	0.93	0.22	0.23	42,63,110,114	0
2	VSB	A	601	36/36	0.94	0.20	-0.11	34,52,97,104	0
4	CL	B	603	1/1	0.98	0.15	-2.14	78,78,78,78	0
4	CL	A	603	1/1	0.85	0.28	-	73,73,73,73	0
3	CA	A	602	1/1	0.53	0.20	-	83,83,83,83	0
3	CA	B	602	1/1	0.42	0.40	-	89,89,89,89	0

6.5 Other polymers [i](#)

There are no such residues in this entry.