



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:23 pm GMT

PDB ID : 1YBW
Title : Protease domain of HGFA with no inhibitor
Authors : Shia, S.; Stamos, J.; Kirchhofer, D.; Fan, B.; Wu, J.; Corpuz, R.T.; Santell, L.; Lazarus, R.A.; Eigenbrot, C.
Deposited on : 2004-12-21
Resolution : 2.70 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	:	1.7.2 (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

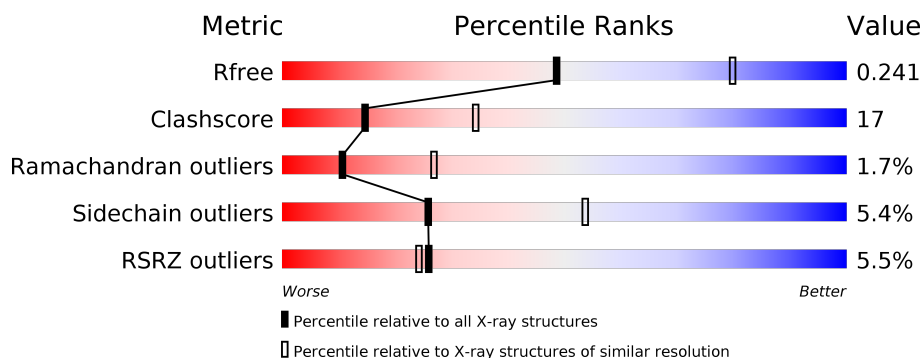
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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	283	<div> <div>2%</div> <div> <div></div> <div>58%</div> <div>27%</div> <div>••</div> <div>13%</div> </div> </div>
1	B	283	<div> <div>7%</div> <div> <div></div> <div>59%</div> <div>26%</div> <div>••</div> <div>13%</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
2	NAG	A	741	-	-	-	X

2 Entry composition [i](#)

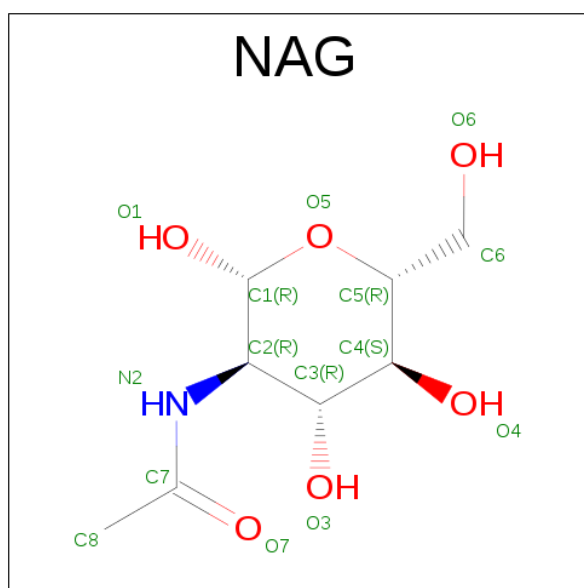
There are 3 unique types of molecules in this entry. The entry contains 3906 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Hepatocyte growth factor activator precursor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	246	Total	C	N	O	S	14	1	0
			1884	1191	331	347	15			
1	B	246	Total	C	N	O	S	17	1	0
			1884	1191	331	347	15			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

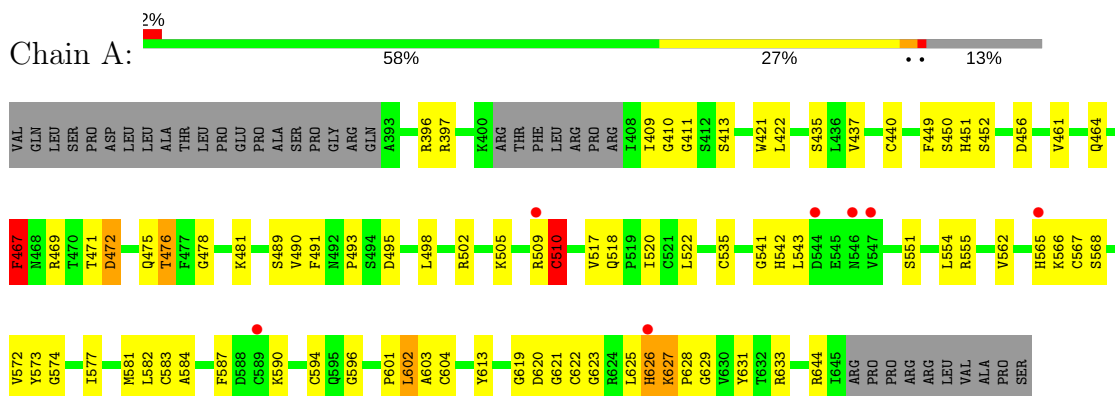
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	57	Total 57	O 57	1	0
3	B	53	Total 53	O 53	1	0

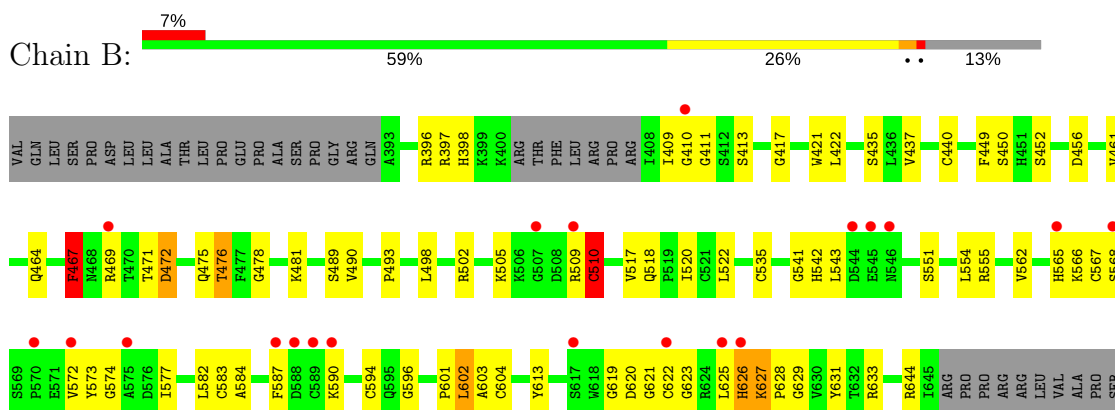
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Hepatocyte growth factor activator precursor



- Molecule 1: Hepatocyte growth factor activator precursor



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.43Å 76.43Å 72.15Å 90.00° 107.77° 90.00°	Depositor
Resolution (Å)	47.94 – 2.70 47.94 – 2.69	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.94-2.70) 99.0 (47.94-2.69)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 2.69Å)	Xtriage
Refinement program	CNX 2000.1	Depositor
R, R_{free}	0.213 , 0.256 0.220 , 0.241	Depositor DCC
R_{free} test set	436 reflections (3.00%)	DCC
Wilson B-factor (Å ²)	31.1	Xtriage
Anisotropy	0.567	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 53.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3906	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.33 % 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 8.8907e-03. 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, 0.333 respectively for untwinned datasets, and 0.375, 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: NAG

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	0.43	0/1938	0.70	3/2634 (0.1%)
1	B	0.42	0/1938	0.71	3/2634 (0.1%)
All	All	0.43	0/3876	0.70	6/5268 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	602	LEU	N-CA-C	-6.47	93.52	111.00
1	A	602	LEU	N-CA-C	-6.46	93.55	111.00
1	A	467	PHE	N-CA-C	5.39	125.55	111.00
1	B	467	PHE	N-CA-C	5.34	125.42	111.00
1	B	478	GLY	N-CA-C	-5.18	100.15	113.10
1	A	478	GLY	N-CA-C	-5.10	100.35	113.10

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	1884	0	1808	61	0
1	B	1884	0	1808	58	0
2	A	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	14	0	13	0	0
3	A	57	0	0	5	0
3	B	53	0	0	5	0
All	All	3906	0	3642	118	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 17.

All (118) 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:471:THR:HG22	1:A:472:ASP:H	1.15	1.12
1:B:471:THR:HG22	1:B:472:ASP:H	1.16	1.06
1:B:567[B]:CYS:SG	1:B:583:CYS:HB2	2.01	1.00
1:A:567[B]:CYS:SG	1:A:583:CYS:HB2	2.01	0.99
1:B:440:CYS:HG	1:B:510:CYS:HG	0.88	0.87
1:A:440:CYS:HG	1:A:510:CYS:HG	0.96	0.87
1:B:509:ARG:O	1:B:510:CYS:HB2	1.79	0.82
1:A:509:ARG:O	1:A:510:CYS:HB2	1.78	0.81
1:A:471:THR:HG22	1:A:472:ASP:N	1.96	0.80
1:B:409:ILE:HG23	1:B:594:CYS:SG	2.22	0.80
1:A:409:ILE:HG23	1:A:594:CYS:SG	2.22	0.80
1:B:471:THR:HG22	1:B:472:ASP:N	1.98	0.78
1:B:627:LYS:HB2	1:B:628:PRO:HD2	1.67	0.77
1:A:627:LYS:HB2	1:A:628:PRO:HD2	1.67	0.76
1:A:535:CYS:HG	1:A:604:CYS:HG	0.74	0.74
1:A:471:THR:CG2	1:A:472:ASP:H	1.96	0.71
1:A:567[B]:CYS:SG	1:A:583:CYS:CB	2.79	0.70
1:B:567[B]:CYS:HB3	1:B:583:CYS:SG	2.33	0.69
1:B:435:SER:OG	1:B:601:PRO:HB3	1.93	0.68
1:A:435:SER:OG	1:A:601:PRO:HB3	1.94	0.68
1:B:567[B]:CYS:SG	1:B:583:CYS:CB	2.80	0.67
1:A:567[B]:CYS:HB3	1:A:583:CYS:SG	2.37	0.64
1:A:498:LEU:HD23	1:A:498:LEU:C	2.18	0.63
1:B:490:VAL:O	1:B:493:PRO:HG3	1.99	0.63
1:A:644:ARG:NE	3:A:36:HOH:O	2.31	0.63
1:B:498:LEU:HD23	1:B:498:LEU:C	2.20	0.62
1:B:502:ARG:HH22	1:B:505:LYS:HE3	1.64	0.61
1:A:490:VAL:O	1:A:493:PRO:HG3	2.00	0.61
1:A:502:ARG:HH22	1:A:505:LYS:HE3	1.64	0.60
1:B:644:ARG:NE	3:B:136:HOH:O	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:577:ILE:HG23	1:B:577:ILE:O	2.02	0.60
1:A:577:ILE:O	1:A:577:ILE:HG23	2.02	0.59
1:B:409:ILE:HD11	1:B:623:GLY:HA3	1.85	0.58
1:B:471:THR:CG2	1:B:472:ASP:H	1.97	0.58
1:A:409:ILE:HD11	1:A:623:GLY:HA3	1.86	0.57
1:B:476:THR:HG23	3:B:160:HOH:O	2.04	0.56
1:B:587:PHE:HA	1:B:626:HIS:O	2.06	0.56
1:A:587:PHE:HA	1:A:626:HIS:O	2.05	0.55
1:A:421:TRP:CG	1:A:520:ILE:HB	2.41	0.55
1:A:475:GLN:NE2	1:A:517:VAL:HG21	2.22	0.54
1:A:396:ARG:O	1:A:518:GLN:NE2	2.40	0.54
1:B:421:TRP:CG	1:B:520:ILE:HB	2.41	0.54
1:A:440:CYS:CB	1:A:510:CYS:HG	2.21	0.54
1:B:396:ARG:O	1:B:518:GLN:NE2	2.40	0.54
1:B:489:SER:HA	3:B:139:HOH:O	2.07	0.53
1:B:475:GLN:NE2	1:B:517:VAL:HG21	2.23	0.53
1:B:464:GLN:NE2	1:B:467:PHE:O	2.42	0.53
1:B:461:VAL:HG13	1:B:476:THR:HG22	1.89	0.53
1:A:464:GLN:NE2	1:A:467:PHE:O	2.42	0.52
1:A:476:THR:HG23	3:A:60:HOH:O	2.09	0.52
1:A:567[B]:CYS:SG	1:A:583:CYS:SG	3.08	0.52
1:B:568:SER:O	1:B:574:GLY:HA2	2.10	0.52
1:A:567[A]:CYS:SG	1:A:577:ILE:HG12	2.50	0.52
1:A:461:VAL:HG13	1:A:476:THR:HG22	1.90	0.52
1:B:567[A]:CYS:SG	1:B:577:ILE:HG12	2.50	0.51
1:A:568:SER:O	1:A:574:GLY:HA2	2.11	0.50
1:B:411:GLY:O	1:B:555:ARG:NH1	2.45	0.49
1:B:440:CYS:CB	1:B:510:CYS:HG	2.24	0.49
1:B:461:VAL:HG22	1:B:476:THR:HB	1.95	0.48
1:B:422:LEU:HD21	1:B:554:LEU:HD11	1.95	0.48
1:B:409:ILE:CD1	1:B:623:GLY:HA3	2.43	0.48
1:B:627:LYS:HB2	1:B:628:PRO:CD	2.42	0.48
1:A:422:LEU:HD21	1:A:554:LEU:HD11	1.96	0.48
1:A:411:GLY:O	1:A:555:ARG:NH1	2.46	0.47
1:B:422:LEU:CD2	1:B:554:LEU:HD11	2.45	0.47
1:B:572:VAL:HG12	1:B:573:TYR:CD2	2.49	0.47
1:A:627:LYS:HB2	1:A:628:PRO:CD	2.42	0.47
1:A:489:SER:HA	3:A:39:HOH:O	2.14	0.47
1:B:449:PHE:O	1:B:452:SER:N	2.45	0.47
1:B:565:HIS:N	3:B:156:HOH:O	2.44	0.47
1:A:461:VAL:HG22	1:A:476:THR:HB	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:409:ILE:CD1	1:A:623:GLY:HA3	2.44	0.47
1:A:437:VAL:HB	1:A:522:LEU:HD21	1.96	0.47
1:A:572:VAL:HG12	1:A:573:TYR:CD2	2.50	0.47
1:B:594:CYS:HA	1:B:619:GLY:O	2.16	0.46
1:B:437:VAL:HB	1:B:522:LEU:HD21	1.97	0.46
1:A:449:PHE:O	1:A:452:SER:N	2.45	0.46
1:A:565:HIS:N	3:A:56:HOH:O	2.45	0.46
1:A:422:LEU:CD2	1:A:554:LEU:HD11	2.45	0.46
1:A:594:CYS:HA	1:A:619:GLY:O	2.15	0.46
1:A:543:LEU:HD23	1:A:543:LEU:O	2.17	0.45
1:A:469:ARG:HH11	1:A:469:ARG:HG3	1.82	0.45
1:B:543:LEU:HD23	1:B:543:LEU:O	2.16	0.45
1:B:567[B]:CYS:SG	1:B:583:CYS:SG	3.15	0.45
1:B:469:ARG:HH11	1:B:469:ARG:HG3	1.83	0.44
1:B:594:CYS:SG	1:B:622:CYS:HA	2.57	0.44
1:B:440:CYS:HG	1:B:510:CYS:CB	2.29	0.44
1:A:518:GLN:HB2	3:A:37:HOH:O	2.16	0.44
1:A:491:PHE:HA	1:B:398:HIS:CD2	2.53	0.44
1:A:594:CYS:SG	1:A:622:CYS:HA	2.58	0.44
1:A:535:CYS:HB3	1:A:603:ALA:O	2.18	0.44
1:B:535:CYS:HB3	1:B:603:ALA:O	2.18	0.43
1:B:582:LEU:HD23	1:B:633:ARG:HG3	1.99	0.43
1:A:620:ASP:O	1:A:621:GLY:C	2.57	0.43
1:B:417:GLY:HA2	3:B:133:HOH:O	2.19	0.43
1:A:568:SER:HA	1:A:577:ILE:HG21	2.01	0.43
1:A:582:LEU:HD23	1:A:633:ARG:HG3	2.00	0.42
1:B:568:SER:HA	1:B:577:ILE:HG21	2.01	0.42
1:A:562:VAL:HG23	1:A:584:ALA:HA	2.02	0.42
1:B:562:VAL:HG23	1:B:584:ALA:HA	2.02	0.42
1:B:449:PHE:O	1:B:450:SER:C	2.58	0.42
1:A:495:ASP:HA	1:A:581:MET:HE2	2.02	0.41
1:A:449:PHE:O	1:A:450:SER:C	2.57	0.41
1:B:541:GLY:HA2	1:B:596:GLY:HA3	2.03	0.41
1:B:604:CYS:SG	1:B:613:TYR:CB	3.08	0.41
1:A:451:HIS:O	1:A:452:SER:C	2.59	0.41
1:A:566:LYS:O	1:A:572:VAL:HG21	2.20	0.41
1:A:498:LEU:HD23	1:A:498:LEU:O	2.19	0.41
1:A:541:GLY:HA2	1:A:596:GLY:HA3	2.02	0.41
1:A:604:CYS:SG	1:A:613:TYR:CB	3.09	0.41
1:B:583:CYS:HA	1:B:629:GLY:O	2.21	0.41
1:A:583:CYS:HA	1:A:629:GLY:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:566:LYS:O	1:B:572:VAL:HG21	2.20	0.41
1:B:620:ASP:O	1:B:621:GLY:C	2.58	0.41
1:A:631:TYR:CD1	1:A:631:TYR:N	2.89	0.40
1:B:410:GLY:HA3	1:B:590:LYS:O	2.21	0.40
1:B:631:TYR:CD1	1:B:631:TYR:N	2.89	0.40
1:A:410:GLY:HA3	1:A:590:LYS:O	2.21	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	243/283 (86%)	212 (87%)	27 (11%)	4 (2%)	11	28
1	B	243/283 (86%)	213 (88%)	26 (11%)	4 (2%)	11	28
All	All	486/566 (86%)	425 (87%)	53 (11%)	8 (2%)	11	28

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	510	CYS
1	B	510	CYS
1	A	467	PHE
1	B	467	PHE
1	A	625	LEU
1	B	625	LEU
1	B	626	HIS
1	A	626	HIS

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	205/237 (86%)	194 (95%)	11 (5%)	26	54
1	B	205/237 (86%)	194 (95%)	11 (5%)	26	54
All	All	410/474 (86%)	388 (95%)	22 (5%)	26	54

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	397	ARG
1	A	413	SER
1	A	456	ASP
1	A	472	ASP
1	A	476	THR
1	A	481	LYS
1	A	510	CYS
1	A	542	HIS
1	A	551	SER
1	A	602	LEU
1	A	627	LYS
1	B	397	ARG
1	B	413	SER
1	B	456	ASP
1	B	472	ASP
1	B	476	THR
1	B	481	LYS
1	B	510	CYS
1	B	542	HIS
1	B	551	SER
1	B	602	LEU
1	B	627	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	642	ASN
1	B	607	ASN
1	B	642	ASN

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	NAG	A	741	1	14,14,15	0.61	0	15,19,21	0.81	1 (6%)
2	NAG	B	742	1	14,14,15	0.61	0	15,19,21	0.77	1 (6%)

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	NAG	A	741	1	-	0/6/23/26	0/1/1/1
2	NAG	B	742	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	741	NAG	C2-N2-C7	-2.25	119.65	122.94
2	B	742	NAG	C2-N2-C7	-2.12	119.85	122.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	246/283 (86%)	0.20	7 (2%) 53 54	9, 25, 56, 68	5 (2%)
1	B	246/283 (86%)	0.35	20 (8%) 13 10	13, 31, 62, 70	6 (2%)
All	All	492/566 (86%)	0.28	27 (5%) 26 24	9, 28, 61, 70	11 (2%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	570	PRO	5.2
1	A	546	ASN	4.5
1	B	626	HIS	3.7
1	B	589	CYS	3.5
1	B	546	ASN	3.1
1	A	547	VAL	3.0
1	B	509	ARG	2.7
1	A	626	HIS	2.7
1	A	565	HIS	2.7
1	B	544	ASP	2.7
1	A	589	CYS	2.6
1	B	625	LEU	2.6
1	A	509	ARG	2.5
1	B	565	HIS	2.5
1	B	590	LYS	2.5
1	B	568	SER	2.4
1	A	544	ASP	2.4
1	B	588	ASP	2.3
1	B	622	CYS	2.2
1	B	410	GLY	2.2
1	B	587	PHE	2.1
1	B	545	GLU	2.1
1	B	469	ARG	2.0
1	B	507	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	617	SER	2.0
1	B	575	ALA	2.0
1	B	572	VAL	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, 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	NAG	A	741	14/15	0.74	0.31	2.48	51,55,57,57	0
2	NAG	B	742	14/15	0.76	0.29	1.65	52,56,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.