



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:37 pm GMT

PDB ID : 3K2U
Title : Crystal structure of HGFA in complex with the allosteric inhibitory antibody Fab40
Authors : Ganesan, R.; Eigenbrot, C.; Shia, S.
Deposited on : 2009-09-30
Resolution : 2.35 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

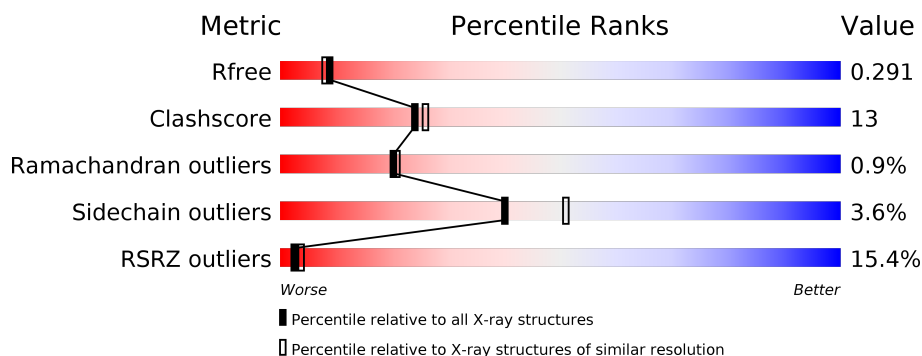
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1522 (2.38-2.34)
Clashscore	112137	1626 (2.38-2.34)
Ramachandran outliers	110173	1605 (2.38-2.34)
Sidechain outliers	110143	1606 (2.38-2.34)
RSRZ outliers	101464	1528 (2.38-2.34)

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	257	<div> <div>18%</div> <div> <div>68%</div> <div>21%</div> <div>• 9%</div> </div> </div>
2	B	35	<div> <div>6%</div> <div> <div>9%</div> <div>6%</div> <div>86%</div> </div> </div>
3	H	225	<div> <div>12%</div> <div> <div>76%</div> <div>18%</div> <div>• •</div> </div> </div>
4	L	214	<div> <div>13%</div> <div> <div>64%</div> <div>31%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5207 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 long chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1780	1133	306	328	13			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	253	ALA	-	EXPRESSION TAG	UNP Q04756
A	254	ALA	-	EXPRESSION TAG	UNP Q04756
A	255	ALA	-	EXPRESSION TAG	UNP Q04756
A	256	HIS	-	EXPRESSION TAG	UNP Q04756
A	257	HIS	-	EXPRESSION TAG	UNP Q04756
A	258	HIS	-	EXPRESSION TAG	UNP Q04756
A	259	HIS	-	EXPRESSION TAG	UNP Q04756
A	260	HIS	-	EXPRESSION TAG	UNP Q04756
A	261	HIS	-	EXPRESSION TAG	UNP Q04756

- Molecule 2 is a protein called Hepatocyte growth factor activator short chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	5	Total	C	N	O	S	0	0	0
			37	20	11	5	1			

- Molecule 3 is a protein called Antibody, Fab fragment, Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	215	Total	C	N	O	S	0	0	0
			1603	1024	265	308	6			

- Molecule 4 is a protein called Antibody, Fab fragment, Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	208	Total	C	N	O	S	0	0	0
			1591	994	267	324	6			

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	2	Total	C	N	O	0	0
			28	16	2	10		

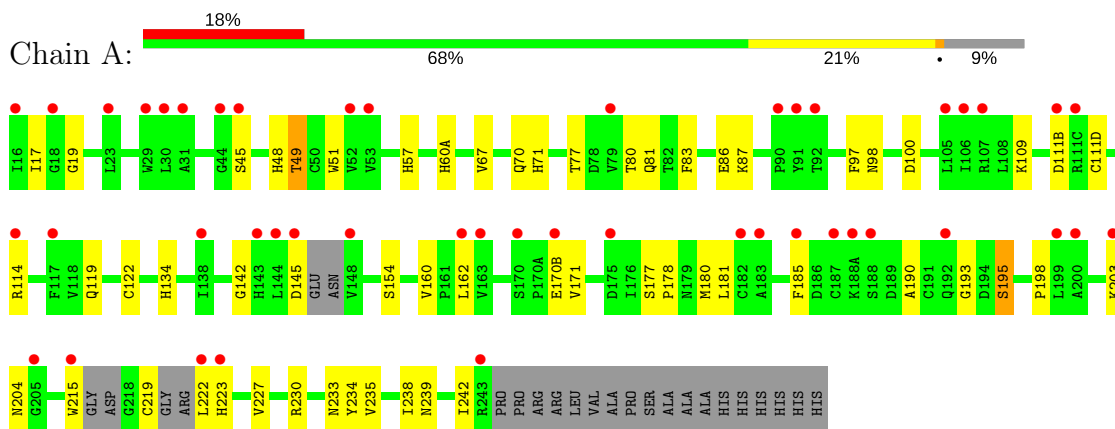
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	72	Total	O	0	0
			72	72		
6	B	2	Total	O	0	0
			2	2		
6	H	54	Total	O	0	0
			54	54		
6	L	40	Total	O	0	0
			40	40		

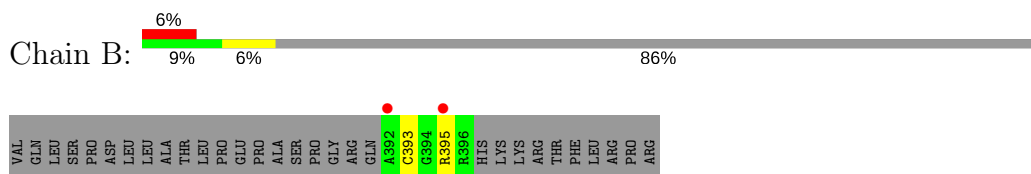
3 Residue-property plots

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.

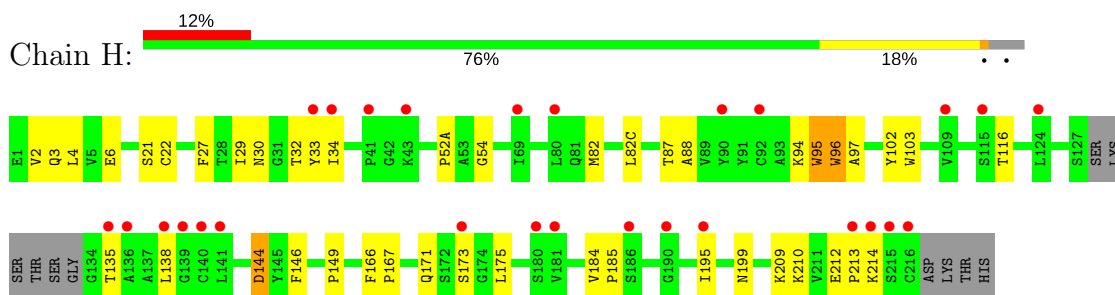
- Molecule 1: Hepatocyte growth factor activator long chain



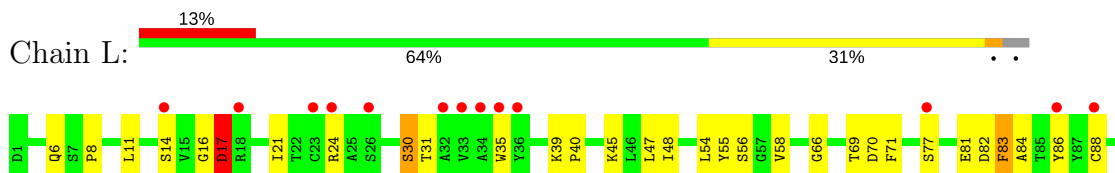
- Molecule 2: Hepatocyte growth factor activator short chain

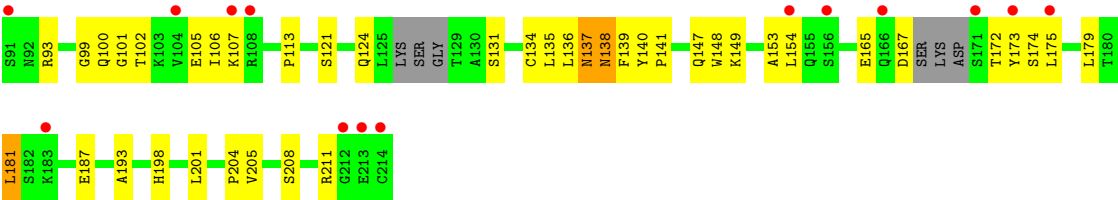


- Molecule 3: Antibody, Fab fragment, Heavy Chain



- Molecule 4: Antibody, Fab fragment, Light Chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	38.94Å 48.93Å 96.03Å 98.10° 95.01° 103.89°	Depositor
Resolution (Å)	19.77 – 2.35 19.77 – 2.35	Depositor EDS
% Data completeness (in resolution range)	94.8 (19.77-2.35) 89.5 (19.77-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.59 (at 2.35Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.239 , 0.290 0.239 , 0.291	Depositor DCC
R_{free} test set	2652 reflections (9.93%)	DCC
Wilson B-factor (Å ²)	49.6	Xtriage
Anisotropy	0.671	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5207	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.52% 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, 0.333 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	0.47	2/1831 (0.1%)	0.60	0/2491
2	B	0.46	0/36	0.70	0/45
3	H	0.45	0/1648	0.64	0/2253
4	L	0.39	0/1623	0.58	0/2202
All	All	0.44	2/5138 (0.0%)	0.61	0/6991

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	17	ILE	C-N	8.15	1.47	1.33
1	A	19	GLY	C-N	5.43	1.46	1.34

There are no bond angle outliers.

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	1780	0	1699	37	0
2	B	37	0	37	1	0
3	H	1603	0	1558	33	0
4	L	1591	0	1540	55	0
5	A	28	0	25	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	72	0	0	4	0
6	B	2	0	0	0	0
6	H	54	0	0	1	0
6	L	40	0	0	2	0
All	All	5207	0	4859	124	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 13.

All (124) 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)
5:A:741:NAG:O6	6:A:312:HOH:O	1.71	1.07
4:L:47:LEU:HD22	4:L:58:VAL:HG13	1.67	0.76
1:A:86:GLU:OE2	1:A:109:LYS:HG2	1.87	0.73
4:L:179:LEU:HG	4:L:181:LEU:HD11	1.75	0.68
4:L:105:GLU:HG2	4:L:106:ILE:N	2.09	0.67
1:A:45:SER:OG	1:A:198:PRO:HB3	1.95	0.67
4:L:6:GLN:HB2	4:L:100:GLN:NE2	2.10	0.67
4:L:11:LEU:O	4:L:11:LEU:HD12	1.95	0.67
3:H:30:ASN:HA	3:H:52(A):PRO:HB2	1.77	0.66
3:H:4:LEU:HD12	3:H:22:CYS:SG	2.36	0.66
1:A:230:ARG:HD2	1:A:233:ASN:ND2	2.11	0.64
1:A:134:HIS:HB3	1:A:162:LEU:HD12	1.80	0.63
3:H:144:ASP:HB3	3:H:175:LEU:HD23	1.82	0.60
3:H:29:ILE:HG22	3:H:34:ILE:HD11	1.83	0.60
4:L:105:GLU:HG2	4:L:106:ILE:H	1.66	0.60
1:A:234:TYR:O	1:A:238:ILE:HG13	2.02	0.60
3:H:214:LYS:HB3	3:H:214:LYS:NZ	2.16	0.59
1:A:233:ASN:HB3	4:L:93:ARG:NH2	2.18	0.59
1:A:114:ARG:HB3	1:A:119:GLN:HE22	1.66	0.59
4:L:140:TYR:CD1	4:L:141:PRO:HA	2.39	0.58
3:H:4:LEU:HD23	3:H:102:TYR:CD1	2.40	0.57
4:L:47:LEU:HD22	4:L:58:VAL:CG1	2.34	0.57
4:L:40:PRO:CB	4:L:165:GLU:HG3	2.36	0.56
4:L:45:LYS:HE3	6:L:241:HOH:O	2.06	0.56
1:A:87:LYS:HB2	3:H:54:GLY:O	2.06	0.55
4:L:81:GLU:C	4:L:83:PHE:H	2.08	0.55
4:L:24:ARG:HG2	4:L:70:ASP:OD1	2.07	0.55
4:L:47:LEU:HA	4:L:58:VAL:HG21	1.88	0.55
3:H:82:MET:HE2	3:H:82(C):LEU:HD21	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:35:TRP:HB2	4:L:48:ILE:HB	1.89	0.54
4:L:16:GLY:HA2	4:L:77:SER:OG	2.08	0.53
5:A:741:NAG:O6	6:A:277:HOH:O	2.15	0.53
1:A:81:GLN:HB3	1:A:83:PHE:HE1	1.73	0.53
1:A:180:MET:CE	1:A:215:TRP:HE1	2.21	0.53
3:H:87:THR:O	3:H:88:ALA:HB2	2.09	0.53
3:H:135:THR:HG22	3:H:185:PRO:HA	1.91	0.52
4:L:6:GLN:HB2	4:L:100:GLN:HE22	1.73	0.52
4:L:187:GLU:HA	4:L:211:ARG:NE	2.24	0.52
3:H:210:LYS:HE3	3:H:212:GLU:HB2	1.90	0.52
4:L:198:HIS:O	4:L:201:LEU:HB2	2.10	0.52
4:L:149:LYS:HA	4:L:153:ALA:O	2.10	0.51
4:L:134:CYS:HB2	4:L:148:TRP:CZ2	2.46	0.51
4:L:11:LEU:C	4:L:11:LEU:HD12	2.32	0.50
3:H:4:LEU:HD23	3:H:102:TYR:HD1	1.76	0.50
4:L:135:LEU:HD11	4:L:137:ASN:HB2	1.94	0.50
4:L:24:ARG:HA	4:L:69:THR:O	2.11	0.50
3:H:6:GLU:HA	3:H:21:SER:O	2.11	0.49
1:A:114:ARG:HB3	1:A:119:GLN:NE2	2.26	0.49
4:L:40:PRO:HB3	4:L:165:GLU:HG3	1.93	0.49
3:H:82:MET:HB3	3:H:82(C):LEU:HD21	1.94	0.49
1:A:60(A):HIS:H	1:A:60(A):HIS:CD2	2.30	0.49
4:L:6:GLN:OE1	4:L:99:GLY:HA3	2.13	0.48
4:L:66:GLY:HA3	4:L:71:PHE:HA	1.95	0.48
1:A:162:LEU:HB3	6:A:301:HOH:O	2.14	0.48
4:L:86:TYR:O	4:L:101:GLY:HA2	2.13	0.48
4:L:39:LYS:HG2	6:L:251:HOH:O	2.14	0.48
3:H:103:TRP:N	3:H:103:TRP:CD1	2.82	0.48
3:H:195:ILE:N	3:H:195:ILE:HD12	2.29	0.47
1:A:180:MET:HE3	1:A:215:TRP:HE1	1.79	0.47
4:L:124:GLN:HE22	4:L:131:SER:N	2.12	0.47
1:A:142:GLY:HA2	1:A:193:GLY:HA3	1.95	0.47
3:H:214:LYS:HZ3	3:H:214:LYS:HB3	1.78	0.47
4:L:167:ASP:HB3	4:L:172:THR:H	1.80	0.47
3:H:116:THR:HA	3:H:146:PHE:O	2.15	0.47
1:A:177:SER:HB2	1:A:178:PRO:HD2	1.95	0.47
4:L:17:ASP:O	4:L:77:SER:HA	2.15	0.47
1:A:171:VAL:HG22	1:A:185:PHE:HE2	1.79	0.46
1:A:49:THR:O	1:A:111(D):CYS:HB2	2.16	0.46
4:L:107:LYS:HA	4:L:140:TYR:OH	2.15	0.46
4:L:124:GLN:HE22	4:L:131:SER:CB	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:212:GLU:HG2	3:H:213:PRO:HD2	1.96	0.46
1:A:86:GLU:O	1:A:87:LYS:HB3	2.17	0.45
1:A:203:LYS:O	1:A:204:ASN:HB2	2.16	0.45
4:L:8:PRO:CG	4:L:11:LEU:HD23	2.46	0.45
1:A:230:ARG:HD2	1:A:233:ASN:HD21	1.80	0.45
3:H:171:GLN:C	3:H:173:SER:H	2.20	0.45
1:A:51:TRP:CD1	1:A:242:ILE:HG23	2.52	0.45
3:H:2:VAL:O	3:H:3:GLN:HG2	2.16	0.45
4:L:83:PHE:O	4:L:84:ALA:HB2	2.17	0.45
4:L:175:LEU:C	4:L:175:LEU:HD23	2.37	0.45
4:L:81:GLU:C	4:L:83:PHE:N	2.70	0.45
3:H:32:THR:OG1	3:H:33:TYR:N	2.50	0.45
1:A:77:THR:OG1	1:A:80:THR:HG23	2.17	0.44
3:H:184:VAL:HB	3:H:185:PRO:HD2	1.99	0.44
1:A:185:PHE:O	1:A:223:HIS:HA	2.18	0.44
4:L:14:SER:O	4:L:17:ASP:HB2	2.18	0.44
1:A:180:MET:HE2	1:A:227:VAL:HG11	1.99	0.44
4:L:21:ILE:HD13	4:L:102:THR:HB	1.99	0.44
4:L:30:SER:HB3	4:L:31:THR:H	1.58	0.44
3:H:27:PHE:HE2	3:H:32:THR:HG21	1.82	0.43
1:A:97:PHE:HB2	3:H:97:ALA:O	2.18	0.43
1:A:181:LEU:C	1:A:181:LEU:HD12	2.38	0.43
4:L:193:ALA:HB2	4:L:208:SER:HB3	2.00	0.43
1:A:190:ALA:O	1:A:219:CYS:HB3	2.19	0.42
4:L:113:PRO:HB3	4:L:139:PHE:HB3	2.01	0.42
3:H:95:TRP:HA	6:H:230:HOH:O	2.19	0.42
4:L:181:LEU:N	4:L:181:LEU:HD12	2.35	0.42
4:L:121:SER:OG	4:L:124:GLN:HB2	2.19	0.42
4:L:124:GLN:HE22	4:L:131:SER:H	1.68	0.42
3:H:195:ILE:HA	3:H:209:LYS:O	2.19	0.42
3:H:195:ILE:HG23	3:H:209:LYS:O	2.19	0.42
4:L:6:GLN:HG2	4:L:88:CYS:SG	2.59	0.42
4:L:154:LEU:HD13	4:L:154:LEU:C	2.40	0.42
4:L:55:TYR:O	4:L:56:SER:C	2.58	0.42
1:A:48:HIS:CE1	1:A:49:THR:HG23	2.55	0.41
1:A:98:ASN:ND2	1:A:100:ASP:HB2	2.35	0.41
1:A:67:VAL:CG1	1:A:70:GLN:HE21	2.32	0.41
4:L:54:LEU:HD11	4:L:58:VAL:CG1	2.50	0.41
3:H:173:SER:O	3:H:175:LEU:HD13	2.19	0.41
1:A:71:HIS:CE1	1:A:154:SER:HB2	2.56	0.41
3:H:95:TRP:CD1	3:H:95:TRP:C	2.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:138:ASN:HA	4:L:173:TYR:O	2.21	0.41
4:L:139:PHE:CD1	4:L:139:PHE:N	2.89	0.41
1:A:235:VAL:HG12	1:A:239:ASN:ND2	2.35	0.41
1:A:122:CYS:SG	2:B:393:CYS:C	2.99	0.41
3:H:212:GLU:HG2	3:H:213:PRO:CD	2.50	0.41
3:H:95:TRP:O	3:H:96:TRP:HB2	2.20	0.41
4:L:40:PRO:HG2	4:L:83:PHE:CE2	2.56	0.41
4:L:201:LEU:HG	4:L:205:VAL:HG23	2.03	0.41
4:L:136:LEU:HD13	4:L:175:LEU:HD22	2.02	0.40
1:A:222:LEU:O	1:A:223:HIS:HB2	2.22	0.40
1:A:180:MET:CE	1:A:227:VAL:HG11	2.52	0.40
3:H:166:PHE:HA	3:H:167:PRO:HD3	1.87	0.40
1:A:195:SER:HB2	6:A:37:HOH: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	225/257 (88%)	204 (91%)	21 (9%)	0	100	100
2	B	3/35 (9%)	3 (100%)	0	0	100	100
3	H	211/225 (94%)	194 (92%)	15 (7%)	2 (1%)	20	21
4	L	202/214 (94%)	181 (90%)	17 (8%)	4 (2%)	9	6
All	All	641/731 (88%)	582 (91%)	53 (8%)	6 (1%)	20	21

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	H	96	TRP
3	H	144	ASP

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Mol	Chain	Res	Type
4	L	17	ASP
4	L	138	ASN
4	L	204	PRO
4	L	82	ASP

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	195/213 (92%)	188 (96%)	7 (4%)	40	51
2	B	3/30 (10%)	2 (67%)	1 (33%)	0	0
3	H	173/182 (95%)	168 (97%)	5 (3%)	48	60
4	L	181/186 (97%)	174 (96%)	7 (4%)	37	47
All	All	552/611 (90%)	532 (96%)	20 (4%)	40	51

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	THR
1	A	57	HIS
1	A	111(B)	ASP
1	A	145	ASP
1	A	160	VAL
1	A	170(B)	GLU
1	A	195	SER
2	B	395	ARG
3	H	94	LYS
3	H	95	TRP
3	H	138	LEU
3	H	149	PRO
3	H	199	ASN
4	L	17	ASP
4	L	30	SER
4	L	83	PHE
4	L	137	ASN

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Mol	Chain	Res	Type
4	L	147	GLN
4	L	174	SER
4	L	181	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	HIS
1	A	60(A)	HIS
1	A	70	GLN
1	A	119	GLN
1	A	143	HIS
1	A	192	GLN
1	A	204	ASN
1	A	233	ASN
3	H	13	GLN
3	H	30	ASN
3	H	199	ASN
4	L	3	GLN
4	L	92	ASN
4	L	124	GLN
4	L	137	ASN
4	L	147	GLN
4	L	189	HIS

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](#)

2 carbohydrates 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	$\# Z > 2$
5	NAG	A	741	1,5	14,14,15	0.55	0	15,19,21	1.10	1 (6%)
5	NAG	A	742	5	14,14,15	0.49	0	15,19,21	2.30	5 (33%)

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
5	NAG	A	741	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	742	5	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	742	NAG	C4-C3-C2	-5.99	102.23	111.02
5	A	741	NAG	O5-C1-C2	-2.75	107.64	111.47
5	A	742	NAG	C2-N2-C7	-2.72	118.98	122.94
5	A	742	NAG	O4-C4-C3	2.07	114.86	110.36
5	A	742	NAG	O4-C4-C5	2.42	115.39	109.28
5	A	742	NAG	C1-O5-C5	4.04	117.73	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	741	NAG	2	0

5.6 Ligand geometry

There are no ligands in this entry.

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	233/257 (90%)	0.96	46 (19%) 1 2	37, 63, 93, 106	1 (0%)
2	B	5/35 (14%)	1.43	2 (40%) 0 0	74, 78, 84, 89	0
3	H	215/225 (95%)	0.68	27 (12%) 4 7	30, 62, 98, 119	1 (0%)
4	L	208/214 (97%)	0.83	27 (12%) 4 6	40, 75, 86, 110	0
All	All	661/731 (90%)	0.83	102 (15%) 2 4	30, 68, 92, 119	2 (0%)

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	216	CYS	10.9
4	L	214	CYS	8.2
1	A	222	LEU	6.5
3	H	69	ILE	6.4
1	A	148	VAL	6.1
1	A	187	CYS	5.8
3	H	190	GLY	5.3
3	H	135	THR	4.7
1	A	223	HIS	4.5
4	L	34	ALA	4.5
3	H	139	GLY	4.5
1	A	143	HIS	4.2
2	B	392	ALA	4.1
1	A	144	LEU	4.0
1	A	192	GLN	4.0
3	H	215	SER	3.9
1	A	79	VAL	3.8
1	A	145	ASP	3.8
3	H	140	CYS	3.4
1	A	106	ILE	3.4
3	H	124	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	31	ALA	3.3
4	L	26	SER	3.3
4	L	35	TRP	3.2
1	A	105	LEU	3.2
2	B	395	ARG	3.2
1	A	111(C)	ARG	3.2
4	L	33	VAL	3.2
3	H	141	LEU	3.2
1	A	111(B)	ASP	3.1
1	A	23	LEU	3.1
3	H	214	LYS	3.1
4	L	104	VAL	3.1
4	L	88	CYS	3.1
1	A	185	PHE	3.1
3	H	90	TYR	3.0
1	A	205	GLY	3.0
1	A	200	ALA	3.0
1	A	18	GLY	3.0
4	L	183	LYS	3.0
3	H	136	ALA	2.9
1	A	53	VAL	2.9
4	L	77	SER	2.9
4	L	36	TYR	2.9
1	A	182	CYS	2.9
3	H	173	SER	2.8
4	L	171	SER	2.8
1	A	29	TRP	2.8
4	L	86	TYR	2.8
1	A	175	ASP	2.7
1	A	199	LEU	2.7
3	H	138	LEU	2.7
1	A	114	ARG	2.7
1	A	44	GLY	2.7
1	A	138	ILE	2.7
1	A	215	TRP	2.6
4	L	166	GLN	2.6
4	L	213	GLU	2.6
3	H	109	VAL	2.6
1	A	243	ARG	2.5
3	H	186	SER	2.4
3	H	92	CYS	2.4
4	L	154	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	170(B)	GLU	2.4
4	L	175	LEU	2.4
3	H	33	TYR	2.3
3	H	213	PRO	2.3
1	A	52	VAL	2.3
1	A	16	ILE	2.3
3	H	34	ILE	2.3
1	A	183	ALA	2.3
4	L	24	ARG	2.3
1	A	90	PRO	2.3
4	L	107	LYS	2.3
3	H	41	PRO	2.2
1	A	162	LEU	2.2
1	A	107	ARG	2.2
4	L	91	SER	2.2
3	H	43	LYS	2.2
1	A	45	SER	2.2
3	H	80	LEU	2.2
3	H	115	SER	2.2
1	A	30	LEU	2.2
1	A	170	SER	2.1
3	H	180	SER	2.1
4	L	108	ARG	2.1
1	A	163	VAL	2.1
3	H	181	VAL	2.1
4	L	14	SER	2.1
4	L	156	SER	2.1
4	L	23	CYS	2.1
1	A	91	TYR	2.1
4	L	212	GLY	2.1
4	L	32	ALA	2.1
1	A	117	PHE	2.1
4	L	18	ARG	2.1
1	A	92	THR	2.1
1	A	203	LYS	2.1
3	H	195	ILE	2.0
4	L	173	TYR	2.0
1	A	188(A)	LYS	2.0
1	A	188	SER	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](#)

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
5	NAG	A	741	14/15	0.78	0.18	0.18	53,64,68,68	0
5	NAG	A	742	14/15	0.75	0.48	-	71,75,84,84	0

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.