



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 10:00 PM BST

PDB ID : 3RRR
Title : Structure of the RSV F protein in the post-fusion conformation
Authors : McLellan, J.S.; Yongping, Y.; Graham, B.S.; Kwong, P.D.
Deposited on : 2011-04-30
Resolution : 2.82 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

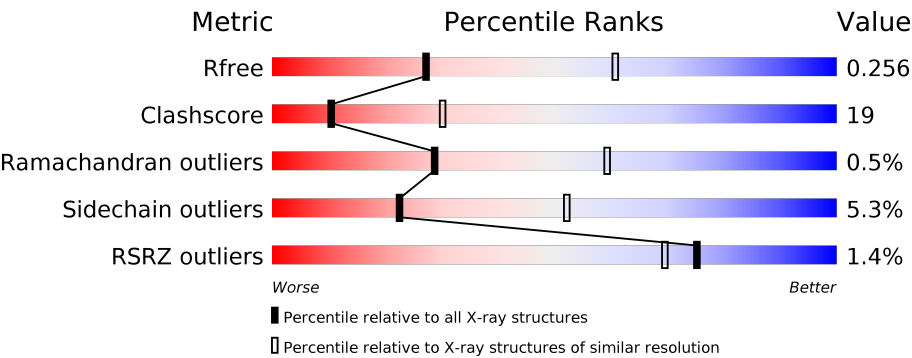
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 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}	130704	3617 (2.84-2.80)
Clashscore	141614	4060 (2.84-2.80)
Ramachandran outliers	138981	3978 (2.84-2.80)
Sidechain outliers	138945	3980 (2.84-2.80)
RSRZ outliers	127900	3552 (2.84-2.80)

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 respectively. 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	84	<div><div>2%</div><div><div></div><div>54%</div><div>31%</div><div>•</div><div>13%</div></div></div>
1	C	84	<div><div></div><div><div>61%</div><div>24%</div><div>•</div><div>14%</div></div></div>
1	E	84	<div><div>%</div><div><div></div><div>64%</div><div>21%</div><div>•</div><div>13%</div></div></div>
1	G	84	<div><div>%</div><div><div></div><div>52%</div><div>33%</div><div></div><div>14%</div></div></div>
1	I	84	<div><div></div><div><div>64%</div><div>19%</div><div>•</div><div>15%</div></div></div>
1	M	84	<div><div>2%</div><div><div></div><div>60%</div><div>26%</div><div></div><div>14%</div></div></div>

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Mol	Chain	Length	Quality of chain
2	B	374	
2	D	374	
2	F	374	
2	H	374	
2	L	374	
2	N	374	

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	NAG	E	770	-	-	-	X
3	NAG	F	800	X	-	-	-
3	NAG	H	800	X	-	-	-
3	NAG	I	770	X	-	-	-
3	NAG	N	800	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 40311 atoms, of which 20196 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 Fusion glycoprotein F0.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	73	Total	C	H	N	O	S	0	0	0
			1167	366	587	95	116	3			
1	C	72	Total	C	H	N	O	S	0	0	0
			1150	361	579	93	114	3			
1	E	73	Total	C	H	N	O	S	0	0	0
			1167	366	587	95	116	3			
1	G	72	Total	C	H	N	O	S	0	0	0
			1150	361	579	93	114	3			
1	I	71	Total	C	H	N	O	S	0	0	0
			1133	356	571	91	112	3			
1	M	72	Total	C	H	N	O	S	0	0	0
			1150	361	579	93	114	3			

- Molecule 2 is a protein called Fusion glycoprotein F0.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	358	Total	C	H	N	O	S	0	0	0
			5563	1744	2798	458	545	18			
2	D	355	Total	C	H	N	O	S	0	0	0
			5532	1735	2783	455	541	18			
2	F	351	Total	C	H	N	O	S	0	0	0
			5483	1720	2757	451	537	18			
2	H	359	Total	C	H	N	O	S	0	0	0
			5582	1750	2809	459	546	18			
2	L	354	Total	C	H	N	O	S	0	0	0
			5527	1733	2780	454	542	18			
2	N	355	Total	C	H	N	O	S	0	0	0
			5539	1737	2787	455	542	18			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	342	TYR	PHE	SEE REMARK 999	UNP Q84850

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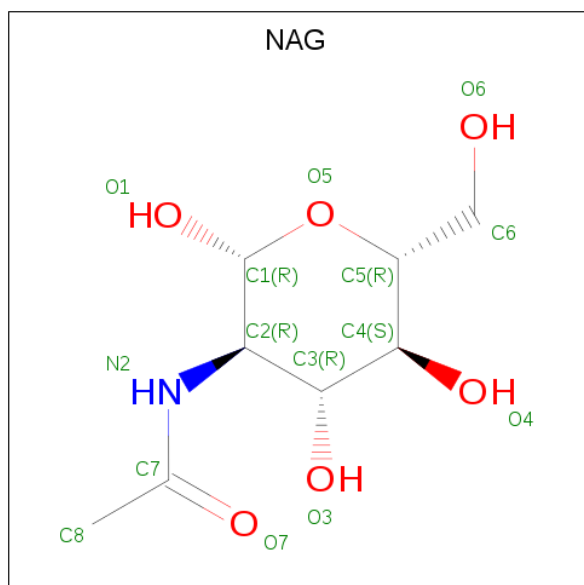
Chain	Residue	Modelled	Actual	Comment	Reference
B	514	GLY	-	expression tag	UNP Q84850
B	515	LEU	-	expression tag	UNP Q84850
B	516	GLU	-	expression tag	UNP Q84850
B	517	VAL	-	expression tag	UNP Q84850
B	518	LEU	-	expression tag	UNP Q84850
B	519	PHE	-	expression tag	UNP Q84850
B	520	GLN	-	expression tag	UNP Q84850
D	342	TYR	PHE	SEE REMARK 999	UNP Q84850
D	514	GLY	-	expression tag	UNP Q84850
D	515	LEU	-	expression tag	UNP Q84850
D	516	GLU	-	expression tag	UNP Q84850
D	517	VAL	-	expression tag	UNP Q84850
D	518	LEU	-	expression tag	UNP Q84850
D	519	PHE	-	expression tag	UNP Q84850
D	520	GLN	-	expression tag	UNP Q84850
F	342	TYR	PHE	SEE REMARK 999	UNP Q84850
F	514	GLY	-	expression tag	UNP Q84850
F	515	LEU	-	expression tag	UNP Q84850
F	516	GLU	-	expression tag	UNP Q84850
F	517	VAL	-	expression tag	UNP Q84850
F	518	LEU	-	expression tag	UNP Q84850
F	519	PHE	-	expression tag	UNP Q84850
F	520	GLN	-	expression tag	UNP Q84850
H	342	TYR	PHE	SEE REMARK 999	UNP Q84850
H	514	GLY	-	expression tag	UNP Q84850
H	515	LEU	-	expression tag	UNP Q84850
H	516	GLU	-	expression tag	UNP Q84850
H	517	VAL	-	expression tag	UNP Q84850
H	518	LEU	-	expression tag	UNP Q84850
H	519	PHE	-	expression tag	UNP Q84850
H	520	GLN	-	expression tag	UNP Q84850
L	342	TYR	PHE	SEE REMARK 999	UNP Q84850
L	514	GLY	-	expression tag	UNP Q84850
L	515	LEU	-	expression tag	UNP Q84850
L	516	GLU	-	expression tag	UNP Q84850
L	517	VAL	-	expression tag	UNP Q84850
L	518	LEU	-	expression tag	UNP Q84850
L	519	PHE	-	expression tag	UNP Q84850
L	520	GLN	-	expression tag	UNP Q84850
N	342	TYR	PHE	SEE REMARK 999	UNP Q84850
N	514	GLY	-	expression tag	UNP Q84850
N	515	LEU	-	expression tag	UNP Q84850

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Chain	Residue	Modelled	Actual	Comment	Reference
N	516	GLU	-	expression tag	UNP Q84850
N	517	VAL	-	expression tag	UNP Q84850
N	518	LEU	-	expression tag	UNP Q84850
N	519	PHE	-	expression tag	UNP Q84850
N	520	GLN	-	expression tag	UNP Q84850

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	F	1	Total	C	N	O	0	0
			14	8	1	5		
3	G	1	Total	C	N	O	0	0
			14	8	1	5		
3	H	1	Total	C	N	O	0	0
			14	8	1	5		

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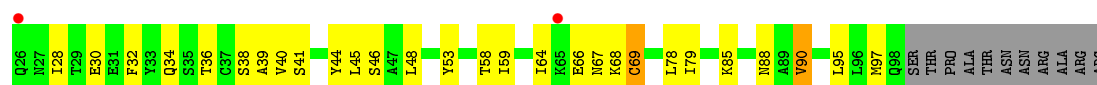
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	I	1	Total	C	N	O	0	0
			14	8	1	5		
3	L	1	Total	C	N	O	0	0
			14	8	1	5		
3	M	1	Total	C	N	O	0	0
			14	8	1	5		
3	N	1	Total	C	N	O	0	0
			14	8	1	5		

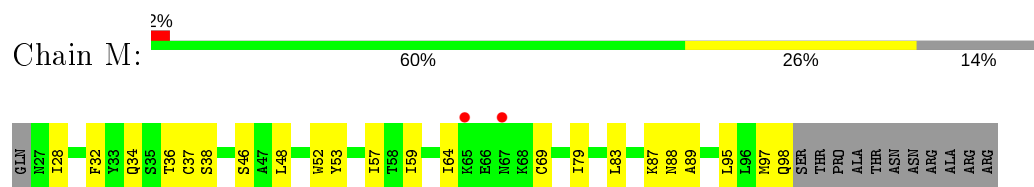
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

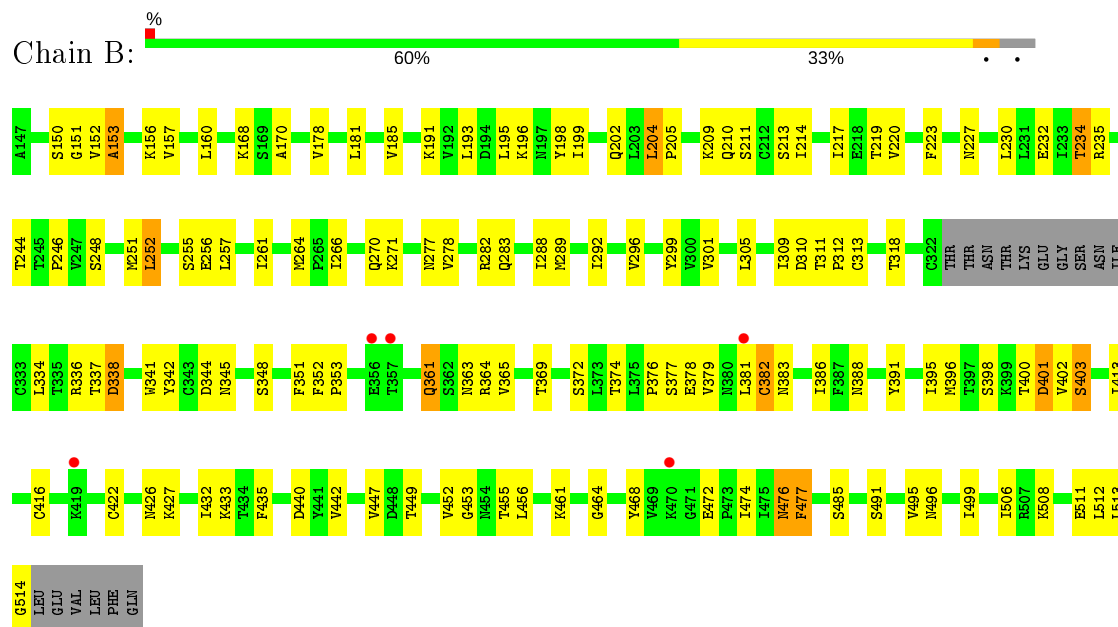
- Molecule 1: Fusion glycoprotein F0



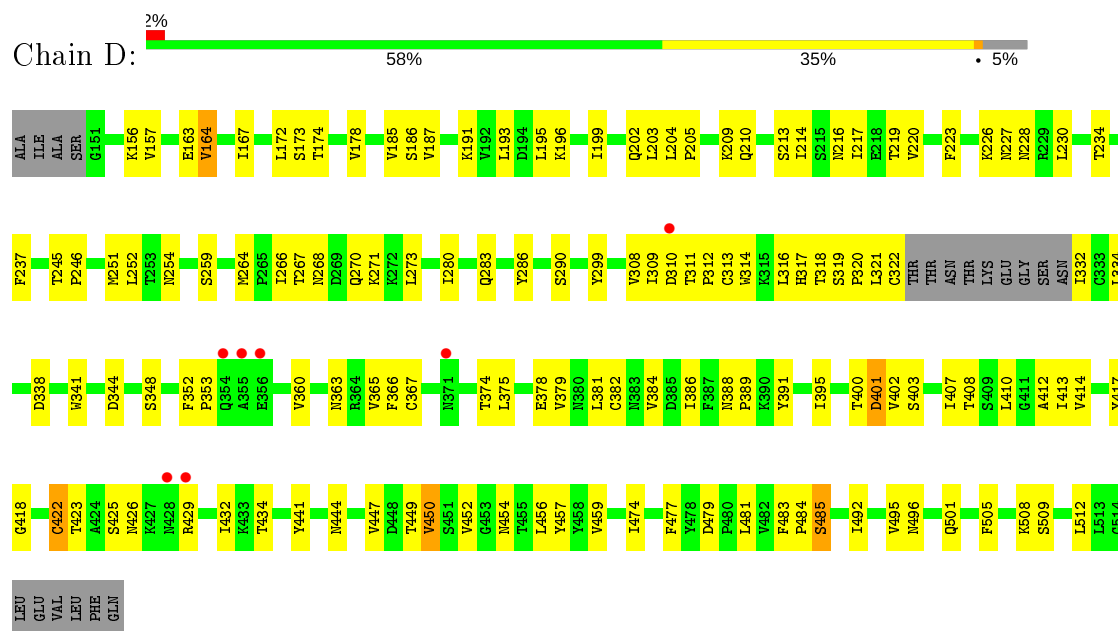
- Molecule 1: Fusion glycoprotein F0



- Molecule 2: Fusion glycoprotein F0

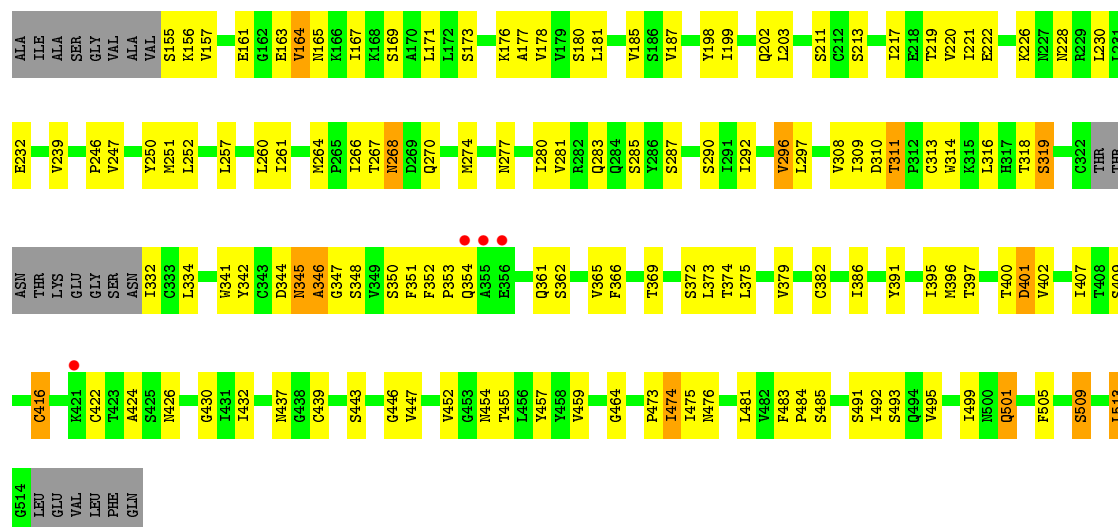


- Molecule 2: Fusion glycoprotein F0

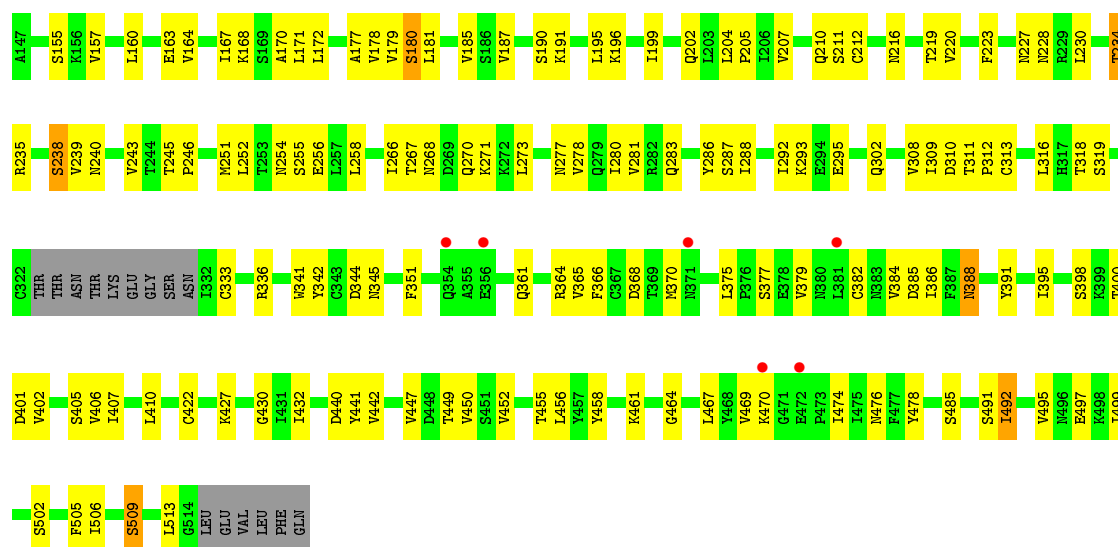


- Molecule 2: Fusion glycoprotein F0

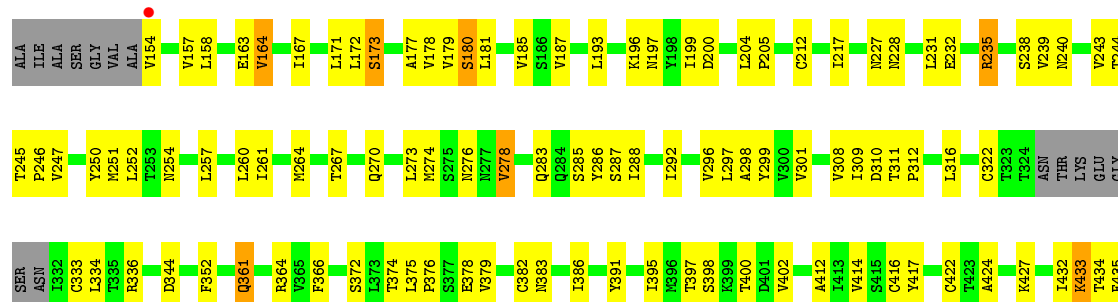




• Molecule 2: Fusion glycoprotein F0

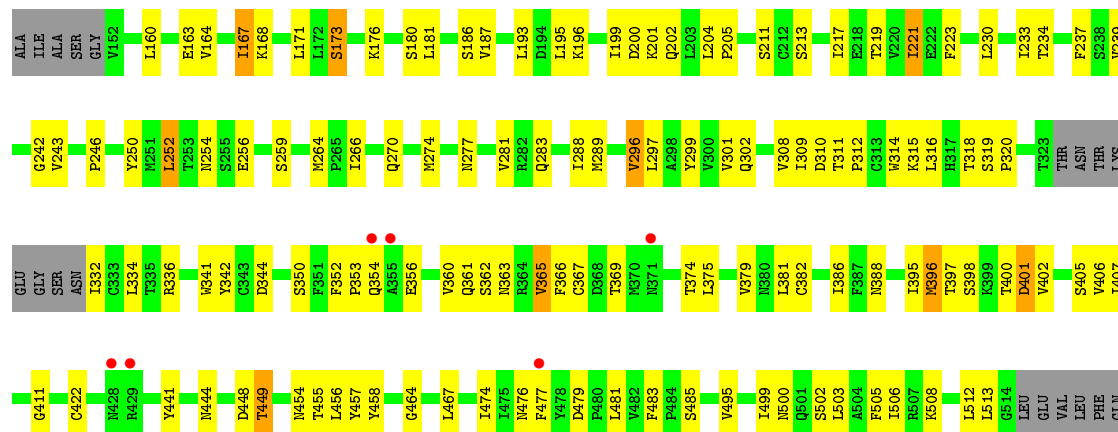


• Molecule 2: Fusion glycoprotein F0





• Molecule 2: Fusion glycoprotein F0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	113.17Å 131.50Å 164.28Å 90.00° 103.17° 90.00°	Depositor
Resolution (Å)	44.21 – 2.82 44.21 – 2.82	Depositor EDS
% Data completeness (in resolution range)	63.7 (44.21-2.82) 63.7 (44.21-2.82)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.221 , 0.262 0.216 , 0.256	Depositor DCC
R_{free} test set	3590 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 28.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	40311	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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	0.27	0/586	0.50	0/789
1	C	0.28	0/577	0.47	0/777
1	E	0.31	0/586	0.45	0/789
1	G	0.27	0/577	0.45	0/777
1	I	0.29	0/568	0.45	0/765
1	M	0.27	0/577	0.46	0/777
2	B	0.29	0/2805	0.47	0/3803
2	D	0.28	0/2789	0.46	0/3781
2	F	0.28	0/2766	0.47	0/3749
2	H	0.28	0/2813	0.47	0/3814
2	L	0.29	0/2787	0.47	0/3779
2	N	0.27	0/2792	0.46	0/3786
All	All	0.28	0/20223	0.46	0/27386

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 [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	580	587	588	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	571	579	579	36	0
1	E	580	587	588	25	0
1	G	571	579	580	33	0
1	I	562	571	572	28	0
1	M	571	579	580	33	0
2	B	2765	2798	2797	128	0
2	D	2749	2783	2782	146	0
2	F	2726	2757	2756	134	0
2	H	2773	2809	2808	145	0
2	L	2747	2780	2779	134	0
2	N	2752	2787	2786	139	0
3	A	14	0	13	0	0
3	B	14	0	13	3	0
3	C	14	0	13	1	0
3	D	14	0	13	0	0
3	E	14	0	13	0	0
3	F	14	0	13	1	0
3	G	14	0	13	1	0
3	H	14	0	13	1	0
3	I	14	0	13	1	0
3	L	14	0	13	0	0
3	M	14	0	13	4	0
3	N	14	0	13	2	0
All	All	20115	20196	20351	753	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 19.

All (753) 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)
2:B:496:ASN:HB3	3:B:800:NAG:H81	1.40	1.01
2:H:308:VAL:HB	2:N:455:THR:HG22	1.60	0.84
2:N:449:THR:HB	2:N:456:LEU:HD11	1.62	0.81
2:N:334:LEU:HB3	2:N:395:ILE:HD11	1.64	0.79
2:H:280:ILE:HD11	2:H:361:GLN:HB2	1.63	0.79
2:H:230:LEU:O	2:H:234:THR:HG23	1.84	0.77
2:F:270:GLN:HG2	2:F:309:ILE:HD12	1.67	0.76
2:L:270:GLN:HG2	2:L:309:ILE:HD12	1.67	0.76
2:N:375:LEU:HB3	2:N:379:VAL:HG11	1.68	0.75
2:L:164:VAL:HA	2:L:167:ILE:HG22	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:495:VAL:CG2	2:D:187:VAL:HG12	2.18	0.74
2:D:308:VAL:HB	2:F:455:THR:HG22	1.70	0.74
2:F:334:LEU:HB3	2:F:395:ILE:HD11	1.70	0.74
1:A:66:GLU:HA	1:A:79:ILE:HG21	1.70	0.73
1:G:28:ILE:HG22	2:H:410:LEU:HD11	1.71	0.73
1:A:46:SER:HB3	2:B:313:CYS:SG	2.29	0.72
1:C:46:SER:HB3	2:D:313:CYS:SG	2.29	0.72
2:H:246:PRO:HB3	2:H:283:GLN:HA	1.70	0.72
1:I:57:ILE:HD11	2:L:252:LEU:HD13	1.73	0.71
2:N:311:THR:HG23	2:N:312:PRO:HD2	1.72	0.71
2:L:333:CYS:HB2	2:L:398:SER:HB3	1.73	0.71
2:L:264:MET:HE3	2:L:274:MET:SD	2.31	0.70
2:F:246:PRO:HB3	2:F:283:GLN:HA	1.72	0.70
2:B:261:ILE:HG12	2:B:264:MET:HE1	1.73	0.70
2:L:334:LEU:HB3	2:L:395:ILE:HD11	1.73	0.70
1:C:36:THR:HG21	2:D:382:CYS:O	1.92	0.70
2:L:386:ILE:HG21	2:L:395:ILE:HD12	1.73	0.69
2:D:477:PHE:CZ	2:D:479:ASP:HB2	2.28	0.69
1:G:38:SER:HB3	2:H:318:THR:HG22	1.75	0.69
2:H:191:LYS:HE3	2:L:491:SER:CB	2.23	0.69
2:N:246:PRO:HB3	2:N:283:GLN:HA	1.74	0.69
2:D:311:THR:HG23	2:D:344:ASP:HB2	1.75	0.68
2:B:230:LEU:O	2:B:234:THR:HG23	1.94	0.68
2:B:246:PRO:HB3	2:B:283:GLN:HA	1.74	0.68
2:D:449:THR:HB	2:D:456:LEU:HD11	1.76	0.68
2:N:264:MET:HE1	2:N:274:MET:SD	2.34	0.68
2:L:311:THR:HG21	2:L:344:ASP:O	1.93	0.67
2:N:270:GLN:HG2	2:N:309:ILE:HD12	1.76	0.67
2:N:318:THR:O	2:N:406:VAL:HG21	1.93	0.67
1:A:45:LEU:O	2:B:364:ARG:HD3	1.94	0.67
2:B:181:LEU:HD23	2:F:181:LEU:HD23	1.75	0.67
1:C:57:ILE:HD11	2:D:252:LEU:HD13	1.77	0.66
1:E:46:SER:HB3	2:F:313:CYS:SG	2.35	0.66
2:H:345:ASN:OD1	2:N:455:THR:HG21	1.94	0.66
2:N:315:LYS:HD2	2:N:341:TRP:CE2	2.30	0.66
2:L:232:GLU:HG3	2:L:250:TYR:CE1	2.30	0.66
2:B:453:GLY:HA3	2:F:346:ALA:HB1	1.78	0.66
1:E:36:THR:HG21	2:F:382:CYS:O	1.95	0.66
2:H:311:THR:HG23	2:H:312:PRO:HD2	1.76	0.66
2:L:177:ALA:HB2	2:N:505:PHE:HB2	1.76	0.66
2:N:252:LEU:CD2	2:N:256:GLU:HB2	2.25	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:270:GLN:HG2	2:B:309:ILE:HD12	1.76	0.65
2:B:334:LEU:HD22	2:B:395:ILE:HD13	1.77	0.65
2:D:246:PRO:HB3	2:D:283:GLN:HA	1.79	0.65
2:H:270:GLN:HG2	2:H:309:ILE:HD12	1.78	0.65
2:H:449:THR:CG2	2:H:456:LEU:HD11	2.25	0.65
2:H:163:GLU:O	2:H:167:ILE:HG22	1.97	0.65
2:H:185:VAL:HG12	2:N:499:ILE:HD11	1.78	0.65
2:L:164:VAL:HG12	2:L:164:VAL:O	1.97	0.65
1:C:57:ILE:CD1	2:D:252:LEU:HD13	2.27	0.65
2:L:288:ILE:N	2:L:288:ILE:HD12	2.11	0.65
2:L:455:THR:HG22	2:N:308:VAL:HB	1.79	0.65
1:I:36:THR:HG22	2:L:386:ILE:HG13	1.77	0.65
2:H:449:THR:HG23	2:H:456:LEU:HD11	1.80	0.64
1:I:36:THR:HG22	2:L:386:ILE:CG1	2.28	0.64
1:M:64:ILE:HG12	1:M:83:LEU:HD21	1.80	0.64
2:D:341:TRP:CH2	2:D:365:VAL:HG21	2.32	0.64
2:F:314:TRP:CE2	2:F:342:TYR:HB2	2.32	0.64
1:G:70:ASN:OD1	3:G:770:NAG:N2	2.31	0.64
1:M:59:ILE:HD12	2:N:233:ILE:HD12	1.80	0.64
2:D:334:LEU:HB3	2:D:395:ILE:HD11	1.80	0.64
2:F:311:THR:HG21	2:F:344:ASP:O	1.98	0.64
2:H:449:THR:CG2	2:H:450:VAL:N	2.59	0.64
2:B:386:ILE:HG21	2:B:395:ILE:HD12	1.77	0.64
1:G:36:THR:HG21	2:H:382:CYS:O	1.97	0.64
1:G:93:LEU:HB3	2:H:292:ILE:HD13	1.80	0.64
2:N:311:THR:HG21	2:N:344:ASP:O	1.98	0.64
2:B:449:THR:HG23	2:B:456:LEU:CD1	2.27	0.64
2:H:280:ILE:HD11	2:H:361:GLN:CB	2.27	0.64
1:I:46:SER:OG	2:L:311:THR:HB	1.98	0.63
2:N:242:GLY:HA2	2:N:289:MET:HE2	1.79	0.63
2:D:505:PHE:HB3	2:F:173:SER:O	1.98	0.63
1:E:37:CYS:SG	2:F:319:SER:HB3	2.38	0.63
2:F:341:TRP:CZ3	2:F:365:VAL:HG21	2.34	0.63
2:H:469:VAL:HB	1:I:59:ILE:HG12	1.80	0.63
2:H:308:VAL:CB	2:N:455:THR:HG22	2.29	0.63
2:D:270:GLN:HG2	2:D:309:ILE:HD12	1.79	0.63
1:A:36:THR:HG21	2:B:382:CYS:O	1.99	0.63
1:I:57:ILE:CD1	2:L:252:LEU:HD13	2.28	0.63
1:M:69:CYS:HA	3:M:770:NAG:H62	1.81	0.63
2:H:499:ILE:HD11	2:L:185:VAL:HG12	1.81	0.62
2:D:375:LEU:HB3	2:D:379:VAL:HG11	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:770:NAG:H3	3:M:770:NAG:O7	2.00	0.62
1:M:83:LEU:HB3	1:M:87:LYS:HE2	1.80	0.62
2:H:199:ILE:HD11	2:L:199:ILE:HD11	1.82	0.62
1:C:64:ILE:HD12	1:C:79:ILE:HG23	1.81	0.61
2:H:235:ARG:O	2:H:239:VAL:HG23	2.00	0.61
2:L:261:ILE:HA	2:L:264:MET:HE2	1.80	0.61
1:M:37:CYS:SG	2:N:319:SER:HB3	2.40	0.61
2:D:401:ASP:OD1	2:D:417:TYR:C	2.38	0.61
2:H:333:CYS:HB2	2:H:398:SER:HB3	1.81	0.61
2:N:311:THR:HG23	2:N:344:ASP:HB2	1.82	0.61
2:L:474:ILE:HB	1:M:64:ILE:HG22	1.83	0.61
2:L:364:ARG:HG3	2:L:366:PHE:CE2	2.35	0.61
1:E:36:THR:HG22	2:F:386:ILE:HG13	1.83	0.61
1:E:90:VAL:HG22	2:F:292:ILE:HD11	1.82	0.60
2:L:477:PHE:CZ	2:L:479:ASP:HB2	2.37	0.60
2:L:442:VAL:CG1	2:L:447:VAL:HB	2.31	0.60
2:B:499:ILE:HD11	2:F:185:VAL:HG12	1.83	0.60
2:N:341:TRP:CH2	2:N:365:VAL:HG21	2.37	0.60
2:B:376:PRO:HB2	2:B:378:GLU:OE1	2.01	0.59
2:H:199:ILE:O	2:H:204:LEU:HB2	2.02	0.59
2:B:168:LYS:HA	2:F:167:ILE:HD11	1.82	0.59
2:D:156:LYS:HB3	2:F:157:VAL:HG11	1.85	0.59
2:F:407:ILE:HD11	2:F:457:TYR:HB3	1.84	0.59
1:G:64:ILE:HD12	1:G:79:ILE:HG23	1.84	0.59
2:D:407:ILE:HD11	2:D:457:TYR:HB3	1.85	0.59
2:H:168:LYS:HA	2:L:167:ILE:HD11	1.84	0.59
1:C:64:ILE:HG22	2:F:474:ILE:CG1	2.31	0.59
2:B:506:ILE:HD11	2:F:178:VAL:HG11	1.83	0.59
2:D:223:PHE:O	2:D:227:ASN:HB2	2.03	0.59
1:M:64:ILE:CG1	1:M:83:LEU:HD21	2.33	0.59
2:B:449:THR:HG23	2:B:456:LEU:HD11	1.84	0.59
1:E:79:ILE:HD13	2:F:220:VAL:HG23	1.85	0.59
2:H:181:LEU:HD21	2:N:181:LEU:HG	1.85	0.59
2:N:386:ILE:HG21	2:N:395:ILE:HD12	1.83	0.59
1:E:97:MET:HE1	2:F:290:SER:O	2.02	0.58
2:F:232:GLU:HG3	2:F:250:TYR:CE1	2.38	0.58
2:N:332:ILE:O	2:N:332:ILE:HG13	2.03	0.58
2:B:449:THR:CG2	2:B:456:LEU:HD11	2.34	0.58
2:H:375:LEU:HD13	2:H:379:VAL:HG11	1.86	0.58
2:L:217:ILE:HG21	2:N:217:ILE:HG21	1.85	0.58
2:N:167:ILE:O	2:N:167:ILE:HG12	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:199:ILE:HD11	2:F:199:ILE:HD11	1.85	0.58
2:B:495:VAL:HG23	2:D:187:VAL:HG12	1.84	0.58
2:B:181:LEU:HD23	2:F:181:LEU:CD2	2.34	0.58
2:L:472:GLU:HB2	2:L:473:PRO:HD2	1.85	0.58
2:B:386:ILE:CG2	2:B:395:ILE:HD12	2.33	0.58
2:D:512:LEU:HD23	2:F:169:SER:HB3	1.85	0.58
2:L:386:ILE:CG2	2:L:395:ILE:HD12	2.34	0.57
1:A:64:ILE:HG22	2:D:474:ILE:HD11	1.86	0.57
1:M:48:LEU:CD2	2:N:367:CYS:HB2	2.34	0.57
2:B:251:MET:HG3	2:B:299:TYR:CE1	2.39	0.57
1:C:79:ILE:HD11	2:D:219:THR:HB	1.86	0.57
2:L:246:PRO:HB3	2:L:283:GLN:HA	1.87	0.57
2:L:442:VAL:HG11	2:L:447:VAL:HG11	1.86	0.57
2:H:280:ILE:HD12	2:H:366:PHE:CE2	2.39	0.57
2:B:341:TRP:CZ3	2:B:365:VAL:HG21	2.40	0.57
1:I:93:LEU:HD11	2:L:238:SER:OG	2.05	0.57
2:N:264:MET:CE	2:N:274:MET:SD	2.92	0.57
2:L:386:ILE:HG21	2:L:395:ILE:CD1	2.35	0.56
2:H:461:LYS:HG2	1:I:52:TRP:HB2	1.86	0.56
2:L:476:ASN:HB3	2:N:219:THR:OG1	2.05	0.56
2:H:273:LEU:HD23	2:H:309:ILE:HD13	1.88	0.56
2:F:277:ASN:O	2:F:281:VAL:HG23	2.05	0.56
2:B:477:PHE:C	2:B:477:PHE:CD2	2.78	0.56
2:H:311:THR:CG2	2:H:344:ASP:HB2	2.36	0.56
2:N:477:PHE:CZ	2:N:479:ASP:HB2	2.39	0.56
2:F:230:LEU:C	2:F:230:LEU:HD23	2.25	0.56
2:L:251:MET:HG3	2:L:299:TYR:CE1	2.41	0.56
2:H:455:THR:HG22	2:L:308:VAL:CG2	2.36	0.56
1:C:30:GLU:HG2	2:D:441:TYR:OH	2.06	0.56
2:H:196:LYS:CE	2:N:485:SER:HB2	2.36	0.56
2:H:318:THR:HG21	2:H:336:ARG:HB2	1.88	0.56
2:L:247:VAL:HG23	2:L:285:SER:HB2	1.88	0.56
1:A:32:PHE:CE2	1:A:34:GLN:HG2	2.41	0.55
1:G:79:ILE:HD11	2:H:219:THR:HB	1.88	0.55
2:H:449:THR:HG22	2:H:450:VAL:N	2.21	0.55
1:E:36:THR:HG22	2:F:386:ILE:CG1	2.37	0.55
2:H:375:LEU:HD23	2:N:396:MET:HE1	1.87	0.55
2:B:278:VAL:O	2:B:282:ARG:HG3	2.06	0.55
1:G:48:LEU:HD12	2:H:308:VAL:CG1	2.36	0.55
2:H:187:VAL:HB	2:L:495:VAL:HG22	1.88	0.55
2:H:384:VAL:CG2	2:H:385:ASP:N	2.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:489:ASP:O	2:L:493:SER:HB2	2.07	0.55
1:C:36:THR:HG22	2:D:386:ILE:HG12	1.88	0.55
2:F:261:ILE:HG12	2:F:274:MET:HE2	1.89	0.55
1:A:78:LEU:HB3	2:B:220:VAL:HG21	1.89	0.55
2:B:181:LEU:O	2:B:185:VAL:HG23	2.07	0.55
2:D:322:CYS:SG	2:D:417:TYR:CD2	2.99	0.55
1:C:48:LEU:CD2	2:D:367:CYS:HB2	2.37	0.55
2:H:375:LEU:HD23	2:N:396:MET:CE	2.37	0.55
2:L:273:LEU:HD23	2:L:309:ILE:HD13	1.88	0.55
2:H:196:LYS:HG2	2:N:483:PHE:CE2	2.42	0.55
2:B:198:TYR:O	2:B:202:GLN:HB2	2.07	0.54
1:A:28:ILE:HD11	2:B:363:ASN:HA	1.87	0.54
2:H:316:LEU:HD11	2:H:336:ARG:HH12	1.72	0.54
2:H:455:THR:HG22	2:L:308:VAL:HG21	1.88	0.54
2:H:311:THR:HG21	2:H:344:ASP:O	2.07	0.54
2:L:247:VAL:HG22	2:L:286:TYR:O	2.07	0.54
2:B:386:ILE:HG21	2:B:395:ILE:CD1	2.36	0.54
2:D:495:VAL:CG2	2:F:187:VAL:HG12	2.36	0.54
1:E:48:LEU:HD12	2:F:345:ASN:HD21	1.72	0.54
2:N:204:LEU:N	2:N:205:PRO:CD	2.70	0.54
2:N:309:ILE:HG22	2:N:310:ASP:CG	2.27	0.54
2:B:352:PHE:CD1	2:B:372:SER:HB3	2.43	0.54
2:D:196:LYS:CE	2:F:485:SER:HB2	2.37	0.54
2:H:311:THR:HG23	2:H:344:ASP:HB2	1.89	0.54
1:A:38:SER:CB	2:B:318:THR:HG22	2.38	0.54
2:D:204:LEU:N	2:D:205:PRO:CD	2.71	0.54
2:D:264:MET:HE3	2:D:266:ILE:CD1	2.38	0.54
1:G:57:ILE:HG12	2:N:467:LEU:HB2	1.90	0.54
2:L:171:LEU:HD13	2:N:171:LEU:HD13	1.89	0.54
2:B:491:SER:HB2	2:D:191:LYS:HE2	1.89	0.54
1:A:85:LYS:HG3	2:D:228:ASN:ND2	2.23	0.54
2:H:204:LEU:N	2:H:205:PRO:CD	2.71	0.54
2:H:308:VAL:HG22	2:H:309:ILE:N	2.22	0.54
1:A:38:SER:HB3	2:B:318:THR:HG22	1.90	0.54
2:B:455:THR:HG22	2:F:308:VAL:CG2	2.37	0.54
1:G:90:VAL:HG22	2:H:292:ILE:HD11	1.89	0.54
2:B:152:VAL:O	2:B:153:ALA:C	2.46	0.54
2:F:309:ILE:CG2	2:F:310:ASP:N	2.72	0.53
2:F:316:LEU:HD22	2:F:318:THR:HG23	1.90	0.53
2:L:412:ALA:HB2	2:L:459:VAL:HG22	1.90	0.53
2:B:513:LEU:HD11	2:F:171:LEU:HD23	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:452:VAL:O	2:F:346:ALA:CB	2.57	0.53
2:B:513:LEU:N	2:B:514:GLY:HA2	2.24	0.53
2:D:374:THR:HG22	2:F:402:VAL:HG21	1.91	0.53
2:H:164:VAL:HA	2:H:167:ILE:HG23	1.90	0.53
2:N:405:SER:HB2	2:N:457:TYR:CE2	2.42	0.53
2:D:210:GLN:O	2:D:214:ILE:HG13	2.08	0.53
2:D:317:HIS:CG	2:D:408:THR:HG22	2.43	0.53
2:F:164:VAL:HA	2:F:167:ILE:HG22	1.91	0.53
2:F:348:SER:HB3	2:F:375:LEU:O	2.09	0.53
2:F:509:SER:O	2:F:513:LEU:HG	2.09	0.53
2:H:497:GLU:HA	3:H:800:NAG:H83	1.90	0.53
2:L:311:THR:HG23	2:L:344:ASP:HB2	1.91	0.53
2:N:266:ILE:HD12	2:N:270:GLN:HB2	1.91	0.53
1:G:64:ILE:CG2	2:N:474:ILE:HD11	2.38	0.53
2:H:160:LEU:HD23	2:N:160:LEU:HD23	1.91	0.53
1:M:48:LEU:HD23	2:N:367:CYS:HB2	1.89	0.53
2:N:309:ILE:HG22	2:N:310:ASP:N	2.22	0.53
2:L:164:VAL:HA	2:L:167:ILE:CG2	2.37	0.52
2:L:433:LYS:NZ	2:L:435:PHE:CE1	2.77	0.52
2:H:476:ASN:HD21	1:I:67:ASN:HB2	1.74	0.52
2:B:232:GLU:O	2:B:235:ARG:HB3	2.08	0.52
2:D:316:LEU:HD21	2:D:318:THR:CG2	2.39	0.52
3:I:770:NAG:O7	3:I:770:NAG:H3	2.08	0.52
2:L:416:CYS:O	2:L:417:TYR:CD1	2.62	0.52
1:C:32:PHE:CG	2:D:441:TYR:HB2	2.44	0.52
1:C:64:ILE:HG22	2:F:474:ILE:HD11	1.91	0.52
2:D:217:ILE:HG23	2:F:221:ILE:HD11	1.92	0.52
2:N:213:SER:O	2:N:217:ILE:HG13	2.10	0.52
1:G:46:SER:OG	2:H:311:THR:HB	2.09	0.52
2:L:402:VAL:HG21	2:N:374:THR:HG22	1.92	0.52
2:B:442:VAL:CG1	2:B:447:VAL:HB	2.39	0.52
2:B:496:ASN:CB	3:B:800:NAG:H81	2.28	0.52
2:D:199:ILE:O	2:D:204:LEU:HB2	2.09	0.52
2:L:361:GLN:HG3	2:L:361:GLN:O	2.10	0.52
1:A:64:ILE:HD12	1:A:79:ILE:HG23	1.90	0.52
1:A:97:MET:SD	2:B:292:ILE:HG22	2.50	0.52
2:D:174:THR:O	2:D:178:VAL:HG23	2.09	0.52
2:H:168:LYS:CA	2:L:167:ILE:HD11	2.39	0.52
2:L:235:ARG:O	2:L:239:VAL:HG23	2.10	0.52
2:N:407:ILE:HD11	2:N:457:TYR:HB3	1.91	0.52
1:A:95:LEU:HB3	2:D:254:ASN:ND2	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:342:TYR:CE1	2:F:351:PHE:CD1	2.98	0.52
2:D:213:SER:O	2:D:217:ILE:HG13	2.11	0.51
2:D:311:THR:HG23	2:D:312:PRO:HD2	1.92	0.51
1:M:83:LEU:O	1:M:87:LYS:HG3	2.09	0.51
2:D:204:LEU:HD21	2:F:481:LEU:HB2	1.92	0.51
1:G:32:PHE:CG	2:H:441:TYR:HB2	2.44	0.51
1:E:83:LEU:HB3	1:E:87:LYS:HE2	1.92	0.51
2:F:361:GLN:NE2	2:F:362:SER:HB2	2.25	0.51
2:L:200:ASP:O	2:L:205:PRO:HD3	2.11	0.51
1:M:37:CYS:SG	1:M:37:CYS:O	2.67	0.51
2:N:411:GLY:HA2	2:N:444:ASN:N	2.26	0.51
1:M:32:PHE:CD1	2:N:441:TYR:HB2	2.45	0.51
2:D:360:VAL:HG13	2:D:360:VAL:O	2.11	0.51
2:H:400:THR:HG22	2:H:401:ASP:N	2.26	0.51
1:M:32:PHE:CG	2:N:441:TYR:HB2	2.46	0.51
2:B:400:THR:HG22	2:B:401:ASP:N	2.26	0.51
2:B:491:SER:CB	2:D:191:LYS:HE2	2.40	0.51
2:D:217:ILE:HG23	2:F:221:ILE:CD1	2.41	0.51
2:H:164:VAL:HA	2:H:167:ILE:CG2	2.39	0.51
1:I:64:ILE:CG1	1:I:83:LEU:HD21	2.41	0.51
2:L:379:VAL:HG12	2:L:391:TYR:CZ	2.46	0.51
2:L:322:CYS:SG	2:L:417:TYR:CD2	3.04	0.51
1:A:36:THR:HB	2:B:336:ARG:HD2	1.93	0.51
2:B:461:LYS:HG2	1:E:52:TRP:HB2	1.93	0.51
2:D:378:GLU:OE2	2:F:400:THR:HG21	2.11	0.51
1:A:68:LYS:O	1:A:69:CYS:HB3	2.10	0.50
2:H:375:LEU:HB3	2:H:379:VAL:HG11	1.93	0.50
2:N:500:ASN:HD22	3:N:800:NAG:C7	2.24	0.50
1:G:52:TRP:CE3	2:H:302:GLN:HG2	2.47	0.50
2:L:376:PRO:HB2	2:L:378:GLU:OE1	2.12	0.50
2:B:248:SER:CB	2:F:239:VAL:HA	2.41	0.50
2:D:441:TYR:CD2	2:D:441:TYR:C	2.85	0.50
2:D:216:ASN:O	2:D:220:VAL:HG23	2.12	0.50
2:N:164:VAL:HA	2:N:167:ILE:HG22	1.92	0.50
2:B:374:THR:HG22	2:D:402:VAL:CG2	2.40	0.50
2:N:252:LEU:HD23	2:N:256:GLU:HB2	1.93	0.50
2:L:180:SER:HB3	2:N:502:SER:HB2	1.94	0.50
2:L:196:LYS:HD3	2:L:196:LYS:C	2.31	0.50
2:N:400:THR:HG22	2:N:401:ASP:H	1.76	0.50
2:D:425:SER:O	2:D:447:VAL:HG13	2.12	0.50
2:D:412:ALA:HB2	2:D:459:VAL:HG13	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:196:LYS:HD3	2:H:196:LYS:C	2.31	0.50
2:L:451:SER:HB2	2:L:456:LEU:HD12	1.93	0.50
1:E:97:MET:C	1:E:98:GLN:HG3	2.31	0.50
2:F:247:VAL:HG23	2:F:285:SER:HB2	1.94	0.50
2:B:170:ALA:CB	2:F:513:LEU:HD21	2.41	0.50
2:H:277:ASN:O	2:H:281:VAL:HG23	2.12	0.50
2:D:264:MET:HE3	2:D:266:ILE:HD13	1.94	0.50
2:L:157:VAL:O	2:L:157:VAL:HG12	2.12	0.50
2:B:210:GLN:O	2:B:214:ILE:HG13	2.12	0.49
2:B:338:ASP:HB2	2:B:342:TYR:OH	2.12	0.49
1:A:95:LEU:HD13	2:D:254:ASN:HD22	1.75	0.49
3:F:800:NAG:O7	3:F:800:NAG:H3	2.11	0.49
2:H:506:ILE:HD11	2:L:178:VAL:HG11	1.94	0.49
2:L:228:ASN:O	2:L:232:GLU:HG2	2.12	0.49
2:D:163:GLU:OE2	2:D:163:GLU:HA	2.12	0.49
2:D:196:LYS:HD3	2:D:196:LYS:C	2.32	0.49
1:G:51:GLY:HA3	2:N:458:TYR:HB2	1.93	0.49
1:G:95:LEU:HD22	2:N:254:ASN:ND2	2.27	0.49
2:D:163:GLU:O	2:D:167:ILE:HG22	2.12	0.49
1:C:46:SER:OG	2:D:311:THR:HB	2.12	0.49
1:E:79:ILE:HD13	2:F:220:VAL:CG2	2.43	0.49
2:B:217:ILE:HG21	2:F:217:ILE:HD13	1.95	0.49
2:H:368:ASP:OD1	2:H:370:MET:HB2	2.13	0.49
2:H:455:THR:HG22	2:L:308:VAL:HB	1.94	0.49
2:N:199:ILE:HG22	2:N:200:ASP:N	2.27	0.49
2:N:309:ILE:CG2	2:N:310:ASP:N	2.75	0.49
2:H:311:THR:CG2	2:H:312:PRO:HD2	2.42	0.49
2:L:477:PHE:CE2	2:L:479:ASP:HB2	2.47	0.49
2:B:309:ILE:CG2	2:B:310:ASP:N	2.76	0.49
2:B:374:THR:HG22	2:D:402:VAL:HG21	1.94	0.49
1:C:64:ILE:HD11	1:C:83:LEU:HG	1.95	0.49
1:M:79:ILE:O	1:M:83:LEU:HG	2.13	0.49
2:D:422:CYS:HB3	2:D:452:VAL:HG22	1.93	0.49
1:I:32:PHE:CG	2:L:441:TYR:HB2	2.47	0.49
2:B:449:THR:CG2	2:B:456:LEU:CD1	2.91	0.49
1:C:32:PHE:CD2	2:D:441:TYR:HB3	2.47	0.49
2:B:351:PHE:CE2	2:B:353:PRO:HB3	2.48	0.49
2:D:217:ILE:HG21	2:F:217:ILE:HG21	1.95	0.49
2:F:222:GLU:O	2:F:226:LYS:HG3	2.13	0.49
2:H:309:ILE:CG2	2:H:310:ASP:N	2.76	0.49
2:L:163:GLU:OE2	2:L:163:GLU:HA	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:267:THR:OG1	2:L:270:GLN:HG3	2.13	0.49
2:N:395:ILE:HG12	2:N:396:MET:N	2.28	0.49
2:B:452:VAL:O	2:F:346:ALA:HB3	2.13	0.48
2:F:296:VAL:HG13	2:F:296:VAL:O	2.12	0.48
2:F:447:VAL:HG12	2:F:459:VAL:HG21	1.95	0.48
2:H:202:GLN:C	2:H:205:PRO:HD2	2.33	0.48
2:H:180:SER:HB3	2:L:502:SER:HB2	1.95	0.48
2:H:171:LEU:HD13	2:N:171:LEU:HD13	1.95	0.48
2:H:495:VAL:CG2	2:N:187:VAL:HG12	2.43	0.48
1:A:95:LEU:HB3	2:D:254:ASN:HD22	1.77	0.48
2:F:252:LEU:HD23	2:F:257:LEU:HD13	1.95	0.48
2:F:422:CYS:SG	2:F:452:VAL:HG22	2.53	0.48
2:N:366:PHE:CD2	2:N:366:PHE:N	2.82	0.48
2:D:311:THR:HG21	2:D:344:ASP:O	2.14	0.48
2:H:485:SER:HB2	2:L:196:LYS:CE	2.44	0.48
2:L:402:VAL:CG2	2:N:374:THR:HG22	2.43	0.48
1:A:48:LEU:HB3	2:B:369:THR:CG2	2.44	0.48
2:F:264:MET:HB2	2:F:264:MET:HE3	1.65	0.48
2:L:247:VAL:HG12	2:L:251:MET:HB3	1.95	0.48
2:B:196:LYS:C	2:B:196:LYS:HD3	2.32	0.48
2:D:280:ILE:HG21	2:D:366:PHE:CD2	2.48	0.48
1:G:61:LEU:HD22	1:G:86:TYR:CZ	2.49	0.48
2:D:273:LEU:HD23	2:D:309:ILE:CD1	2.44	0.48
2:L:375:LEU:HD13	2:L:379:VAL:HG11	1.96	0.48
2:D:319:SER:OG	2:D:320:PRO:HD2	2.13	0.48
2:L:240:ASN:HB3	2:L:243:VAL:O	2.13	0.48
2:N:449:THR:CB	2:N:456:LEU:HD11	2.36	0.48
2:N:476:ASN:OD1	2:N:476:ASN:N	2.47	0.48
2:D:273:LEU:HD23	2:D:309:ILE:HD13	1.95	0.48
2:D:348:SER:HB3	2:D:375:LEU:O	2.14	0.48
2:F:228:ASN:O	2:F:232:GLU:HG2	2.14	0.48
2:F:334:LEU:HD22	2:F:395:ILE:CD1	2.44	0.48
2:L:481:LEU:O	2:N:204:LEU:HD21	2.13	0.48
2:L:492:ILE:HG13	2:N:193:LEU:HD13	1.96	0.48
1:A:36:THR:HG22	2:B:386:ILE:CG1	2.44	0.48
2:D:388:ASN:OD1	2:D:389:PRO:HD2	2.14	0.48
2:B:277:ASN:OD1	2:B:361:GLN:HG2	2.14	0.47
2:N:264:MET:HE3	2:N:266:ILE:HD13	1.94	0.47
2:N:316:LEU:HD21	2:N:318:THR:CG2	2.44	0.47
2:L:173:SER:O	2:N:505:PHE:HB3	2.13	0.47
2:H:172:LEU:HD23	2:N:513:LEU:HD12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:332:ILE:O	2:D:332:ILE:HG13	2.14	0.47
2:D:492:ILE:O	2:D:496:ASN:ND2	2.46	0.47
2:B:311:THR:HG23	2:B:344:ASP:HB2	1.96	0.47
1:C:48:LEU:HD23	2:D:367:CYS:HB2	1.96	0.47
2:D:483:PHE:CD1	2:D:484:PRO:HD2	2.49	0.47
2:H:432:ILE:HD11	2:H:447:VAL:HG22	1.95	0.47
1:C:53:TYR:CE1	2:F:464:GLY:HA3	2.49	0.47
2:D:204:LEU:HD23	2:D:204:LEU:O	2.13	0.47
1:C:45:LEU:HD22	2:D:310:ASP:HA	1.97	0.47
1:A:58:THR:HG22	1:A:59:ILE:N	2.29	0.47
1:A:79:ILE:HD11	2:B:219:THR:HB	1.95	0.47
2:L:336:ARG:HH12	2:L:383:ASN:ND2	2.12	0.47
2:D:401:ASP:C	2:D:401:ASP:OD2	2.52	0.47
2:D:348:SER:OG	2:F:402:VAL:HG23	2.13	0.47
1:G:28:ILE:CD1	1:G:44:TYR:CE1	2.97	0.47
2:L:187:VAL:HB	2:N:495:VAL:HG22	1.97	0.47
2:L:424:ALA:O	2:L:432:ILE:HG12	2.14	0.47
2:D:316:LEU:CD2	2:D:318:THR:CG2	2.93	0.47
2:L:254:ASN:HD22	1:M:95:LEU:HD13	1.79	0.47
1:C:32:PHE:CD2	2:D:441:TYR:CB	2.97	0.47
2:D:187:VAL:O	2:D:191:LYS:HG2	2.14	0.47
2:D:251:MET:HG3	2:D:299:TYR:CE1	2.50	0.47
1:I:69:CYS:O	2:L:212:CYS:SG	2.72	0.47
2:L:297:LEU:HD12	2:L:298:ALA:N	2.30	0.47
1:M:46:SER:OG	2:N:311:THR:HB	2.14	0.47
2:B:156:LYS:O	2:B:160:LEU:HD13	2.15	0.47
2:H:240:ASN:HB3	2:H:243:VAL:O	2.15	0.47
1:C:63:ASN:O	2:F:473:PRO:HA	2.15	0.47
1:E:48:LEU:HB2	2:F:308:VAL:HG12	1.97	0.47
2:L:451:SER:HB2	2:L:456:LEU:CD1	2.45	0.47
1:G:57:ILE:HD11	2:H:252:LEU:HD13	1.97	0.47
2:H:407:ILE:HD13	2:H:458:TYR:O	2.14	0.47
1:A:53:TYR:CD2	2:B:264:MET:HG2	2.50	0.46
2:B:337:THR:CG2	2:B:396:MET:HB2	2.44	0.46
2:B:363:ASN:C	2:B:363:ASN:OD1	2.53	0.46
2:B:336:ARG:HH12	2:B:383:ASN:HD22	1.62	0.46
2:D:230:LEU:O	2:D:234:THR:HG23	2.15	0.46
1:G:44:TYR:HB2	2:H:313:CYS:HB2	1.97	0.46
2:D:410:LEU:HA	2:D:444:ASN:ND2	2.29	0.46
2:H:312:PRO:HG2	2:H:344:ASP:OD2	2.14	0.46
2:H:505:PHE:CD1	2:N:176:LYS:HB2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:53:TYR:CZ	2:N:464:GLY:HA3	2.51	0.46
1:A:46:SER:CB	2:B:313:CYS:SG	3.02	0.46
2:B:476:ASN:HB3	2:F:219:THR:OG1	2.15	0.46
2:D:378:GLU:HG3	2:F:400:THR:OG1	2.15	0.46
2:H:336:ARG:HD2	2:H:386:ILE:HD11	1.96	0.46
2:H:386:ILE:HG21	2:H:395:ILE:HD12	1.96	0.46
2:L:185:VAL:O	2:L:185:VAL:CG1	2.64	0.46
2:L:252:LEU:HD22	2:L:301:VAL:HG11	1.97	0.46
2:H:177:ALA:HB2	2:L:505:PHE:HB2	1.96	0.46
2:D:164:VAL:HA	2:D:167:ILE:CG2	2.46	0.46
2:D:508:LYS:O	2:D:512:LEU:HD13	2.15	0.46
2:H:252:LEU:HD12	2:H:256:GLU:HB2	1.98	0.46
2:L:167:ILE:HG23	2:N:167:ILE:CD1	2.45	0.46
2:N:361:GLN:NE2	2:N:362:SER:HB2	2.30	0.46
2:B:168:LYS:CA	2:F:167:ILE:HD11	2.46	0.46
1:I:53:TYR:CD1	2:L:264:MET:HG2	2.50	0.46
1:I:93:LEU:HB2	2:L:292:ILE:HD13	1.97	0.46
1:I:36:THR:HG21	2:L:382:CYS:O	2.16	0.46
2:B:193:LEU:HD13	2:D:492:ILE:HG13	1.97	0.46
2:B:311:THR:HG23	2:B:312:PRO:HD2	1.97	0.46
2:N:360:VAL:HG22	2:N:361:GLN:N	2.31	0.46
2:B:464:GLY:HA3	1:E:53:TYR:CZ	2.50	0.46
2:F:309:ILE:HG22	2:F:310:ASP:N	2.30	0.46
2:N:199:ILE:O	2:N:204:LEU:HB2	2.15	0.46
2:F:345:ASN:O	2:F:347:GLY:N	2.49	0.46
2:L:312:PRO:HG2	2:L:344:ASP:OD2	2.16	0.46
2:L:352:PHE:CD1	2:L:372:SER:HB3	2.51	0.46
2:B:381:LEU:HB2	2:B:388:ASN:ND2	2.31	0.46
2:B:508:LYS:O	2:B:512:LEU:HD13	2.16	0.46
1:C:33:TYR:CD2	1:C:33:TYR:N	2.84	0.46
2:D:316:LEU:HD23	2:D:318:THR:HG23	1.98	0.46
2:H:342:TYR:CZ	2:H:351:PHE:CD1	3.04	0.46
3:C:770:NAG:H3	3:C:770:NAG:O7	2.15	0.45
1:E:26:GLN:CG	1:E:27:ASN:H	2.29	0.45
2:H:181:LEU:HD23	2:L:181:LEU:HD23	1.98	0.45
2:B:426:ASN:OD1	2:B:427:LYS:N	2.49	0.45
2:F:161:GLU:HG2	2:F:161:GLU:O	2.16	0.45
2:H:379:VAL:HG12	2:H:391:TYR:CE2	2.52	0.45
2:H:455:THR:HG22	2:L:308:VAL:CB	2.46	0.45
2:B:248:SER:HB3	2:F:239:VAL:HA	1.97	0.45
2:F:352:PHE:CE1	2:F:372:SER:HB3	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:48:LEU:O	1:G:50:THR:HG23	2.16	0.45
2:D:379:VAL:HG12	2:D:391:TYR:CZ	2.50	0.45
2:D:501:GLN:O	2:D:505:PHE:CD2	2.70	0.45
2:D:505:PHE:CD1	2:F:176:LYS:HB3	2.52	0.45
2:F:264:MET:HE1	2:F:274:MET:SD	2.56	0.45
2:F:351:PHE:HD2	2:F:373:LEU:HD12	1.81	0.45
2:F:475:ILE:HG22	2:F:476:ASN:N	2.31	0.45
2:L:442:VAL:HG11	2:L:447:VAL:CG1	2.46	0.45
2:D:167:ILE:HG12	2:D:167:ILE:O	2.16	0.45
2:D:423:THR:HG22	2:D:434:THR:HA	1.99	0.45
2:F:292:ILE:HB	2:F:297:LEU:HD13	1.99	0.45
2:L:199:ILE:CD1	2:N:199:ILE:HD11	2.47	0.45
2:L:414:VAL:CG2	2:L:450:VAL:HG11	2.46	0.45
2:N:277:ASN:O	2:N:281:VAL:HG23	2.17	0.45
2:D:334:LEU:HD13	2:D:395:ILE:CD1	2.46	0.45
2:H:223:PHE:CE1	2:H:227:ASN:ND2	2.85	0.45
2:N:319:SER:OG	2:N:320:PRO:CD	2.65	0.45
1:M:28:ILE:HD11	2:N:363:ASN:HA	1.97	0.45
2:N:363:ASN:C	2:N:363:ASN:OD1	2.55	0.45
2:B:433:LYS:HD3	2:B:435:PHE:CE1	2.52	0.45
1:C:64:ILE:HD11	1:C:83:LEU:CG	2.47	0.45
1:I:57:ILE:HD12	2:L:301:VAL:HG21	1.98	0.45
2:L:164:VAL:CG1	2:L:164:VAL:O	2.65	0.45
2:F:198:TYR:O	2:F:202:GLN:HB2	2.17	0.45
2:N:288:ILE:O	2:N:299:TYR:HB2	2.17	0.45
2:B:511:GLU:HG2	2:B:511:GLU:O	2.17	0.45
1:G:83:LEU:O	1:G:87:LYS:HG3	2.17	0.45
2:H:467:LEU:HB2	1:I:57:ILE:HG12	1.98	0.45
2:D:203:LEU:HD21	2:F:203:LEU:CD2	2.47	0.44
2:B:402:VAL:HG13	2:F:374:THR:HG22	1.98	0.44
2:F:386:ILE:HG21	2:F:395:ILE:HD12	1.99	0.44
2:H:168:LYS:HB2	2:L:167:ILE:HD12	1.99	0.44
2:H:254:ASN:O	2:H:258:LEU:HD13	2.17	0.44
2:B:223:PHE:CD1	2:B:227:ASN:ND2	2.85	0.44
1:C:64:ILE:HG22	2:F:474:ILE:HG12	1.99	0.44
2:D:163:GLU:HB3	2:F:164:VAL:HG11	1.99	0.44
2:F:416:CYS:O	2:F:437:ASN:HA	2.18	0.44
2:N:500:ASN:ND2	3:N:800:NAG:C7	2.80	0.44
2:B:491:SER:O	2:B:495:VAL:HG23	2.17	0.44
2:D:426:ASN:HB3	2:D:429:ARG:HB2	1.99	0.44
2:B:191:LYS:HE3	2:F:491:SER:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:64:ILE:HD11	1:G:83:LEU:HG	2.00	0.44
2:H:402:VAL:HG13	2:L:374:THR:CG2	2.48	0.44
1:M:36:THR:HG22	2:N:386:ILE:HG13	1.99	0.44
1:M:38:SER:HB3	2:N:318:THR:HG22	1.99	0.44
1:A:44:TYR:HB2	2:B:313:CYS:HB2	1.98	0.44
2:D:341:TRP:CZ3	2:D:365:VAL:HG21	2.52	0.44
2:D:450:VAL:HG23	2:D:459:VAL:CG2	2.47	0.44
1:G:49:ARG:NH1	1:G:52:TRP:CE2	2.86	0.44
2:H:266:ILE:HD12	2:H:270:GLN:HB2	1.99	0.44
1:M:36:THR:HG21	2:N:382:CYS:O	2.17	0.44
1:C:41:SER:O	2:D:314:TRP:HA	2.18	0.44
1:C:64:ILE:HG22	2:F:474:ILE:CD1	2.48	0.44
2:D:374:THR:HG22	2:F:402:VAL:CG2	2.48	0.44
2:F:379:VAL:HG12	2:F:391:TYR:CE2	2.52	0.44
2:F:400:THR:HG22	2:F:401:ASP:N	2.32	0.44
2:L:171:LEU:CD1	2:N:171:LEU:HD13	2.47	0.44
2:F:501:GLN:O	2:F:505:PHE:HD2	1.99	0.44
2:H:384:VAL:HG23	2:H:385:ASP:N	2.32	0.44
2:H:379:VAL:HG12	2:H:391:TYR:CZ	2.53	0.44
1:M:32:PHE:CD2	2:N:441:TYR:HB3	2.52	0.44
2:B:266:ILE:HG13	2:B:271:LYS:HG3	2.00	0.44
2:B:432:ILE:HD11	2:B:447:VAL:HG22	2.00	0.44
2:D:378:GLU:HB2	2:D:391:TYR:HB2	1.99	0.44
2:H:227:ASN:O	2:H:228:ASN:C	2.54	0.44
2:L:461:LYS:HG2	1:M:52:TRP:HB2	1.99	0.44
3:M:770:NAG:C3	3:M:770:NAG:O7	2.65	0.44
2:L:485:SER:HB2	2:N:196:LYS:NZ	2.32	0.44
2:N:296:VAL:HG13	2:N:296:VAL:O	2.18	0.44
2:B:205:PRO:O	2:B:209:LYS:HG3	2.17	0.44
2:B:318:THR:HG21	2:B:336:ARG:HB2	1.98	0.44
2:B:468:TYR:CZ	2:H:470:LYS:HE3	2.53	0.44
2:D:400:THR:HG22	2:D:401:ASP:N	2.33	0.44
2:D:226:LYS:HD2	2:F:474:ILE:HG21	1.99	0.44
1:I:64:ILE:HG13	1:I:83:LEU:HD11	2.00	0.44
2:H:502:SER:HB2	2:N:180:SER:HB3	1.99	0.44
2:D:185:VAL:HG12	2:F:499:ILE:CD1	2.48	0.44
2:D:191:LYS:HA	2:D:191:LYS:HD2	1.85	0.44
2:D:413:ILE:HG22	2:D:414:VAL:N	2.33	0.44
2:F:342:TYR:CE1	2:F:351:PHE:HD1	2.35	0.44
2:B:278:VAL:O	2:B:282:ARG:CG	2.66	0.43
2:B:348:SER:OG	2:D:402:VAL:HG23	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:196:LYS:HE2	2:D:485:SER:HB2	2.00	0.43
2:F:266:ILE:HD12	2:F:270:GLN:HB3	1.99	0.43
2:D:193:LEU:HD22	2:F:492:ILE:HG21	2.00	0.43
2:H:267:THR:HB	2:H:270:GLN:HG3	2.00	0.43
2:H:308:VAL:CG2	2:H:309:ILE:N	2.81	0.43
2:H:492:ILE:HD13	2:H:492:ILE:HA	1.81	0.43
2:H:506:ILE:HG21	2:L:179:VAL:HG23	2.00	0.43
2:L:227:ASN:O	2:L:231:LEU:HG	2.18	0.43
1:M:52:TRP:CE3	2:N:302:GLN:HG2	2.53	0.43
2:B:379:VAL:HG12	2:B:391:TYR:CZ	2.54	0.43
2:B:452:VAL:O	2:B:452:VAL:HG12	2.18	0.43
1:C:75:LYS:NZ	2:H:427:LYS:HD2	2.33	0.43
2:D:309:ILE:CG2	2:D:310:ASP:N	2.80	0.43
1:I:93:LEU:HB2	2:L:292:ILE:CD1	2.48	0.43
2:N:217:ILE:O	2:N:221:ILE:HG13	2.16	0.43
2:H:179:VAL:HG23	2:N:506:ILE:HG21	1.99	0.43
1:C:37:CYS:HB2	2:D:321:LEU:HD13	2.01	0.43
2:D:401:ASP:OD1	2:D:418:GLY:N	2.51	0.43
2:D:505:PHE:HB2	2:F:177:ALA:HB2	1.99	0.43
1:E:57:ILE:HD11	2:F:252:LEU:HD13	1.99	0.43
2:N:163:GLU:HA	2:N:163:GLU:OE2	2.19	0.43
2:N:361:GLN:O	2:N:361:GLN:HG3	2.16	0.43
2:H:308:VAL:CG1	2:N:455:THR:HG22	2.49	0.43
1:A:39:ALA:HB2	2:B:413:ILE:HD11	1.99	0.43
2:B:496:ASN:C	3:B:800:NAG:H81	2.38	0.43
2:F:345:ASN:O	2:F:346:ALA:C	2.56	0.43
2:N:239:VAL:CG1	2:N:239:VAL:O	2.65	0.43
2:N:334:LEU:CD2	2:N:397:THR:HG22	2.49	0.43
2:B:499:ILE:CD1	2:F:185:VAL:HG12	2.47	0.43
1:C:66:GLU:HG3	1:C:79:ILE:HG21	2.01	0.43
1:E:64:ILE:CG1	1:E:83:LEU:HD21	2.48	0.43
2:F:424:ALA:HB1	2:F:447:VAL:HG21	2.01	0.43
1:G:79:ILE:CD1	2:H:219:THR:CG2	2.97	0.43
2:H:207:VAL:HG12	2:H:207:VAL:O	2.19	0.43
2:H:292:ILE:HG23	2:H:292:ILE:O	2.19	0.43
2:L:276:ASN:HA	1:M:98:GLN:HE22	1.82	0.43
2:N:314:TRP:CE2	2:N:342:TYR:HB2	2.53	0.43
2:B:213:SER:HB3	2:D:214:ILE:HG21	2.01	0.43
1:E:48:LEU:HD13	2:F:369:THR:CG2	2.47	0.43
1:I:61:LEU:HD12	2:L:227:ASN:OD1	2.18	0.43
2:L:264:MET:CE	2:L:274:MET:SD	3.04	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:481:LEU:HD12	2:N:481:LEU:HA	1.89	0.43
1:A:58:THR:CG2	1:A:59:ILE:N	2.81	0.43
2:B:252:LEU:HD12	2:B:301:VAL:HG11	2.00	0.43
2:D:317:HIS:CD2	2:D:408:THR:HG22	2.53	0.43
2:D:452:VAL:HG12	2:D:452:VAL:O	2.17	0.43
1:G:28:ILE:HD11	1:G:44:TYR:CE1	2.54	0.43
2:H:210:GLN:NE2	2:N:211:SER:HB2	2.34	0.43
2:L:464:GLY:HA3	1:M:53:TYR:CZ	2.54	0.43
2:D:316:LEU:CD2	2:D:318:THR:HG23	2.49	0.43
2:N:318:THR:HG21	2:N:336:ARG:HB2	2.00	0.43
2:D:319:SER:OG	2:D:320:PRO:CD	2.67	0.43
2:H:266:ILE:HG13	2:H:271:LYS:HG3	2.00	0.43
1:I:40:VAL:HG22	2:L:316:LEU:HD13	2.00	0.43
1:A:48:LEU:HB3	2:B:369:THR:HG23	2.01	0.43
2:D:267:THR:O	2:D:271:LYS:HG3	2.19	0.43
2:F:400:THR:HG22	2:F:401:ASP:H	1.84	0.43
2:F:430:GLY:O	2:F:432:ILE:HG23	2.19	0.43
2:H:492:ILE:HG21	2:L:193:LEU:HD22	2.01	0.43
2:L:442:VAL:HG11	2:L:447:VAL:CB	2.49	0.43
2:L:432:ILE:HD11	2:L:447:VAL:CG2	2.49	0.43
2:N:400:THR:HG22	2:N:401:ASP:N	2.34	0.43
2:H:196:LYS:NZ	2:N:485:SER:HB2	2.33	0.43
2:B:252:LEU:HD23	2:B:256:GLU:HB2	2.01	0.42
2:H:216:ASN:O	2:H:220:VAL:HG23	2.19	0.42
2:H:318:THR:O	2:H:406:VAL:HG21	2.19	0.42
1:M:48:LEU:HB3	2:N:369:THR:HG23	2.01	0.42
1:M:32:PHE:CG	2:N:441:TYR:CB	3.02	0.42
2:H:251:MET:HG2	2:H:287:SER:CB	2.48	0.42
2:B:252:LEU:O	2:B:282:ARG:NH1	2.50	0.42
2:F:334:LEU:CD2	2:F:397:THR:HG22	2.49	0.42
2:H:388:ASN:O	2:H:388:ASN:ND2	2.52	0.42
2:H:513:LEU:CD1	2:L:172:LEU:HD21	2.49	0.42
2:L:379:VAL:HG12	2:L:391:TYR:CE2	2.55	0.42
1:M:59:ILE:HD12	2:N:233:ILE:CD1	2.48	0.42
2:N:201:LYS:O	2:N:205:PRO:HG2	2.18	0.42
2:B:150:SER:HB2	2:B:151:GLY:HA2	2.02	0.42
1:C:66:GLU:HA	2:F:476:ASN:ND2	2.34	0.42
2:D:267:THR:HG22	2:D:268:ASN:N	2.35	0.42
2:D:167:ILE:HD11	2:F:167:ILE:HG23	2.00	0.42
2:F:452:VAL:HG12	2:F:452:VAL:O	2.20	0.42
2:N:508:LYS:O	2:N:512:LEU:HD13	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:ILE:O	2:B:296:VAL:HG23	2.20	0.42
2:L:196:LYS:HD3	2:L:197:ASN:N	2.34	0.42
2:B:199:ILE:HD11	2:D:199:ILE:HD11	2.01	0.42
2:D:509:SER:HB2	2:F:173:SER:HB3	2.01	0.42
2:H:341:TRP:CH2	2:H:365:VAL:HG21	2.54	0.42
1:A:36:THR:HG22	2:B:386:ILE:HG13	2.02	0.42
2:B:204:LEU:N	2:B:205:PRO:CD	2.82	0.42
2:B:402:VAL:CG1	2:B:403:SER:N	2.82	0.42
1:C:95:LEU:O	1:C:96:LEU:HD23	2.19	0.42
2:F:483:PHE:CG	2:F:484:PRO:HD2	2.55	0.42
2:H:179:VAL:HG22	2:N:503:LEU:CD2	2.49	0.42
2:B:264:MET:HE3	2:B:266:ILE:CD1	2.49	0.42
1:C:83:LEU:O	1:C:87:LYS:HG3	2.20	0.42
1:E:37:CYS:SG	1:E:37:CYS:O	2.77	0.42
2:L:157:VAL:CG1	2:L:157:VAL:O	2.67	0.42
2:L:442:VAL:CG1	2:L:447:VAL:CB	2.97	0.42
2:B:337:THR:HG21	2:B:396:MET:HB2	2.01	0.42
1:E:46:SER:OG	2:F:311:THR:HB	2.20	0.42
1:G:91:THR:O	1:G:95:LEU:HD12	2.20	0.42
1:G:93:LEU:CB	2:H:292:ILE:HD13	2.46	0.42
1:I:64:ILE:HD12	1:I:79:ILE:HG23	2.02	0.42
2:H:170:ALA:CB	2:L:513:LEU:HD21	2.50	0.42
2:B:213:SER:O	2:B:217:ILE:HG13	2.20	0.42
1:A:53:TYR:HB2	2:B:305:LEU:HD11	2.02	0.42
1:E:26:GLN:CG	1:E:27:ASN:N	2.83	0.42
2:H:178:VAL:HG12	2:H:178:VAL:O	2.19	0.42
2:H:386:ILE:CG2	2:H:395:ILE:HD12	2.49	0.42
1:I:64:ILE:HG12	1:I:83:LEU:HD21	2.01	0.42
1:M:57:ILE:HD12	2:N:301:VAL:CG2	2.50	0.42
2:N:405:SER:HB2	2:N:457:TYR:CD2	2.55	0.42
1:C:53:TYR:CZ	2:F:464:GLY:HA3	2.55	0.41
2:D:363:ASN:OD1	2:D:363:ASN:C	2.59	0.41
1:E:64:ILE:HG12	1:E:83:LEU:HD21	2.02	0.41
1:G:30:GLU:HG3	1:G:40:VAL:O	2.19	0.41
1:G:57:ILE:HG12	2:N:467:LEU:HD12	2.01	0.41
1:G:71:GLY:HA2	2:H:212:CYS:HB3	2.01	0.41
2:H:430:GLY:O	2:H:432:ILE:HG23	2.19	0.41
1:I:89:ALA:O	1:I:93:LEU:HD13	2.20	0.41
2:N:316:LEU:CD2	2:N:318:THR:HG23	2.49	0.41
1:A:40:VAL:HG12	1:A:41:SER:N	2.34	0.41
2:D:202:GLN:C	2:D:205:PRO:HD2	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:191:LYS:HE3	2:L:491:SER:HB2	1.99	0.41
2:H:464:GLY:HA3	1:I:53:TYR:CZ	2.55	0.41
1:I:53:TYR:CD2	1:I:53:TYR:C	2.94	0.41
2:H:505:PHE:CE1	2:N:176:LYS:HB3	2.56	0.41
2:B:157:VAL:HG22	2:D:157:VAL:HG22	2.03	0.41
2:D:449:THR:CB	2:D:456:LEU:HD11	2.49	0.41
2:F:395:ILE:HG12	2:F:396:MET:N	2.36	0.41
2:L:309:ILE:CG2	2:L:310:ASP:N	2.83	0.41
1:C:79:ILE:CD1	2:D:219:THR:HB	2.48	0.41
2:D:264:MET:HE3	2:D:266:ILE:HD11	2.02	0.41
2:L:154:VAL:HG23	2:L:154:VAL:O	2.19	0.41
2:L:171:LEU:O	2:L:171:LEU:HG	2.20	0.41
2:L:257:LEU:CD2	2:L:278:VAL:HG12	2.50	0.41
1:M:97:MET:HB2	1:M:97:MET:HE2	1.90	0.41
2:D:172:LEU:CD2	2:F:513:LEU:HD12	2.51	0.41
2:B:248:SER:HB2	2:F:239:VAL:HA	2.02	0.41
2:H:243:VAL:HG22	2:H:288:ILE:HG23	2.03	0.41
2:N:252:LEU:HD21	2:N:256:GLU:HB2	2.02	0.41
2:N:256:GLU:O	2:N:259:SER:HB3	2.21	0.41
2:F:260:LEU:O	2:F:264:MET:HG3	2.21	0.41
2:F:280:ILE:HG21	2:F:366:PHE:CD2	2.55	0.41
2:H:167:ILE:O	2:H:167:ILE:CG1	2.69	0.41
2:H:223:PHE:CD1	2:H:227:ASN:ND2	2.88	0.41
2:H:267:THR:HG22	2:H:268:ASN:N	2.35	0.41
2:H:476:ASN:ND2	1:I:67:ASN:HB2	2.35	0.41
2:H:402:VAL:HG13	2:L:374:THR:HG22	2.03	0.41
2:F:163:GLU:O	2:F:165:ASN:N	2.53	0.41
2:D:505:PHE:HD1	2:F:176:LYS:HB2	1.86	0.41
1:E:79:ILE:CD1	2:F:220:VAL:HG23	2.50	0.41
1:E:93:LEU:HB2	2:F:292:ILE:HD13	2.03	0.41
2:F:426:ASN:ND2	2:F:446:GLY:O	2.54	0.41
2:N:230:LEU:O	2:N:234:THR:HG23	2.21	0.41
2:B:257:LEU:HD23	2:B:278:VAL:CG1	2.51	0.41
2:D:449:THR:C	2:D:450:VAL:HG23	2.41	0.41
2:H:273:LEU:CD2	2:H:309:ILE:HD13	2.50	0.41
1:I:55:SER:HB3	2:L:260:LEU:HD13	2.03	0.41
2:L:287:SER:C	2:L:288:ILE:HD12	2.40	0.41
1:M:88:ASN:O	1:M:89:ALA:C	2.59	0.41
2:N:315:LYS:HD2	2:N:341:TRP:NE1	2.36	0.41
1:C:64:ILE:CD1	1:C:79:ILE:HG23	2.51	0.41
2:D:172:LEU:HD23	2:F:513:LEU:HD12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:245:THR:HB	2:D:286:TYR:CD1	2.56	0.41
2:D:381:LEU:O	2:D:384:VAL:HG22	2.21	0.41
2:B:455:THR:HG22	2:F:308:VAL:HG21	2.02	0.41
2:B:296:VAL:O	2:B:296:VAL:HG13	2.21	0.41
2:H:207:VAL:O	2:H:207:VAL:CG1	2.68	0.41
2:H:509:SER:HB2	2:N:173:SER:HB2	2.03	0.41
1:M:69:CYS:CA	3:M:770:NAG:H62	2.49	0.41
2:N:164:VAL:O	2:N:168:LYS:HB2	2.21	0.41
2:N:252:LEU:HD12	2:N:301:VAL:HG11	2.03	0.41
2:N:352:PHE:HA	2:N:353:PRO:HD2	1.95	0.41
1:A:90:VAL:HG22	2:B:292:ILE:HD11	2.01	0.41
2:H:293:LYS:HE2	2:H:295:GLU:OE2	2.21	0.41
2:H:442:VAL:CG1	2:H:447:VAL:HG21	2.51	0.41
2:L:485:SER:HB2	2:N:196:LYS:CE	2.51	0.41
2:N:381:LEU:HB2	2:N:388:ASN:ND2	2.36	0.41
1:A:36:THR:HG22	2:B:386:ILE:HG12	2.02	0.40
1:A:38:SER:HB2	2:B:318:THR:HG22	2.03	0.40
2:D:365:VAL:O	2:D:365:VAL:HG12	2.20	0.40
2:F:267:THR:HG22	2:F:268:ASN:N	2.35	0.40
1:G:40:VAL:CG1	1:G:42:LYS:HG2	2.52	0.40
2:H:245:THR:HB	2:H:286:TYR:CD1	2.56	0.40
2:N:164:VAL:HA	2:N:167:ILE:CG2	2.51	0.40
2:B:178:VAL:HG12	2:B:178:VAL:O	2.21	0.40
2:D:352:PHE:HA	2:D:353:PRO:HD2	1.83	0.40
2:D:477:PHE:CE2	2:D:479:ASP:HB2	2.56	0.40
2:F:495:VAL:O	2:F:499:ILE:HG13	2.21	0.40
2:H:238:SER:HB3	2:N:250:TYR:CE2	2.56	0.40
2:B:204:LEU:CD2	2:D:481:LEU:HB3	2.51	0.40
1:C:71:GLY:C	2:D:209:LYS:HB3	2.41	0.40
2:L:177:ALA:HB3	2:N:506:ILE:HD11	2.03	0.40
2:L:427:LYS:HG2	2:L:448:ASP:OD2	2.22	0.40
2:B:257:LEU:HD23	2:B:278:VAL:HG12	2.04	0.40
2:F:332:ILE:HG13	2:F:332:ILE:O	2.21	0.40
2:N:195:LEU:HD23	2:N:195:LEU:HA	1.89	0.40
1:C:36:THR:CG2	2:D:382:CYS:O	2.66	0.40
2:F:251:MET:HG2	2:F:287:SER:OG	2.21	0.40
2:F:311:THR:HG23	2:F:344:ASP:HB2	2.02	0.40
2:D:185:VAL:HG12	2:F:499:ILE:HD11	2.03	0.40
2:H:405:SER:HB2	2:H:452:VAL:HG21	2.03	0.40
2:H:491:SER:O	2:H:495:VAL:HG23	2.21	0.40
2:L:375:LEU:HB3	2:L:376:PRO:HD2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:444:ASN:HB2	2:L:460:ASN:OD1	2.21	0.40
1:M:64:ILE:HG21	2:N:223:PHE:HB2	2.04	0.40
2:N:233:ILE:CG2	2:N:297:LEU:HD21	2.51	0.40
2:N:319:SER:OG	2:N:320:PRO:HD2	2.20	0.40
2:N:354:GLN:C	2:N:356:GLU:H	2.24	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	71/84 (84%)	65 (92%)	5 (7%)	1 (1%)	11	32
1	C	70/84 (83%)	68 (97%)	2 (3%)	0	100	100
1	E	71/84 (84%)	65 (92%)	5 (7%)	1 (1%)	11	32
1	G	70/84 (83%)	65 (93%)	5 (7%)	0	100	100
1	I	69/84 (82%)	65 (94%)	4 (6%)	0	100	100
1	M	70/84 (83%)	64 (91%)	6 (9%)	0	100	100
2	B	354/374 (95%)	336 (95%)	17 (5%)	1 (0%)	41	70
2	D	351/374 (94%)	326 (93%)	23 (7%)	2 (1%)	25	54
2	F	347/374 (93%)	320 (92%)	23 (7%)	4 (1%)	13	37
2	H	355/374 (95%)	327 (92%)	28 (8%)	0	100	100
2	L	350/374 (94%)	331 (95%)	17 (5%)	2 (1%)	25	54
2	N	351/374 (94%)	328 (93%)	22 (6%)	1 (0%)	41	70
All	All	2529/2748 (92%)	2360 (93%)	157 (6%)	12 (0%)	29	59

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	69	CYS
2	B	153	ALA
2	F	164	VAL
2	F	346	ALA
2	F	353	PRO
2	L	437	ASN
2	D	450	VAL
2	L	164	VAL
2	D	164	VAL
2	F	296	VAL
2	N	296	VAL
1	E	76	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	66/75 (88%)	62 (94%)	4 (6%)	18	46
1	C	65/75 (87%)	64 (98%)	1 (2%)	65	88
1	E	66/75 (88%)	64 (97%)	2 (3%)	41	73
1	G	65/75 (87%)	65 (100%)	0	100	100
1	I	64/75 (85%)	63 (98%)	1 (2%)	62	87
1	M	65/75 (87%)	64 (98%)	1 (2%)	65	88
2	B	329/344 (96%)	304 (92%)	25 (8%)	13	35
2	D	328/344 (95%)	315 (96%)	13 (4%)	31	64
2	F	326/344 (95%)	304 (93%)	22 (7%)	16	41
2	H	330/344 (96%)	310 (94%)	20 (6%)	18	46
2	L	329/344 (96%)	310 (94%)	19 (6%)	20	48
2	N	329/344 (96%)	311 (94%)	18 (6%)	21	50
All	All	2362/2514 (94%)	2236 (95%)	126 (5%)	22	52

All (126) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLU
1	A	67	ASN
1	A	88	ASN
1	A	90	VAL
2	B	195	LEU
2	B	204	LEU
2	B	211	SER
2	B	234	THR
2	B	244	THR
2	B	252	LEU
2	B	255	SER
2	B	288	ILE
2	B	289	MET
2	B	338	ASP
2	B	345	ASN
2	B	361	GLN
2	B	377	SER
2	B	382	CYS
2	B	398	SER
2	B	401	ASP
2	B	403	SER
2	B	416	CYS
2	B	422	CYS
2	B	440	ASP
2	B	472	GLU
2	B	474	ILE
2	B	476	ASN
2	B	477	PHE
2	B	485	SER
1	C	33	TYR
2	D	173	SER
2	D	186	SER
2	D	195	LEU
2	D	237	PHE
2	D	259	SER
2	D	290	SER
2	D	338	ASP
2	D	401	ASP
2	D	403	SER
2	D	422	CYS
2	D	432	ILE
2	D	454	ASN
2	D	485	SER

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Mol	Chain	Res	Type
1	E	26	GLN
1	E	34	GLN
2	F	155	SER
2	F	156	LYS
2	F	180	SER
2	F	211	SER
2	F	213	SER
2	F	268	ASN
2	F	311	THR
2	F	319	SER
2	F	345	ASN
2	F	350	SER
2	F	354	GLN
2	F	401	ASP
2	F	409	SER
2	F	416	CYS
2	F	439	CYS
2	F	443	SER
2	F	454	ASN
2	F	474	ILE
2	F	493	SER
2	F	501	GLN
2	F	509	SER
2	F	513	LEU
2	H	155	SER
2	H	157	VAL
2	H	180	SER
2	H	190	SER
2	H	195	LEU
2	H	211	SER
2	H	234	THR
2	H	238	SER
2	H	255	SER
2	H	278	VAL
2	H	319	SER
2	H	364	ARG
2	H	377	SER
2	H	388	ASN
2	H	422	CYS
2	H	440	ASP
2	H	474	ILE
2	H	478	TYR

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Mol	Chain	Res	Type
2	H	492	ILE
2	H	509	SER
1	I	67	ASN
2	L	158	LEU
2	L	173	SER
2	L	180	SER
2	L	204	LEU
2	L	235	ARG
2	L	244	THR
2	L	245	THR
2	L	278	VAL
2	L	296	VAL
2	L	361	GLN
2	L	397	THR
2	L	400	THR
2	L	422	CYS
2	L	433	LYS
2	L	434	THR
2	L	440	ASP
2	L	449	THR
2	L	487	GLU
2	L	493	SER
1	M	34	GLN
2	N	167	ILE
2	N	173	SER
2	N	186	SER
2	N	202	GLN
2	N	221	ILE
2	N	237	PHE
2	N	243	VAL
2	N	252	LEU
2	N	350	SER
2	N	365	VAL
2	N	396	MET
2	N	398	SER
2	N	401	ASP
2	N	402	VAL
2	N	422	CYS
2	N	448	ASP
2	N	449	THR
2	N	454	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (59) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	34	GLN
1	A	67	ASN
2	B	159	HIS
2	B	240	ASN
2	B	371	ASN
2	B	383	ASN
2	B	388	ASN
2	B	437	ASN
1	C	34	GLN
2	D	228	ASN
2	D	254	ASN
2	D	276	ASN
2	D	383	ASN
2	D	428	ASN
2	D	437	ASN
2	D	444	ASN
2	D	454	ASN
2	D	476	ASN
2	D	496	ASN
1	E	26	GLN
1	E	34	GLN
1	E	67	ASN
2	F	202	GLN
2	F	254	ASN
2	F	270	GLN
2	F	345	ASN
2	F	383	ASN
2	F	437	ASN
2	F	454	ASN
2	F	476	ASN
1	G	26	GLN
1	G	34	GLN
2	H	270	GLN
2	H	276	ASN
2	H	277	ASN
2	H	383	ASN
2	H	428	ASN
2	H	454	ASN
1	I	34	GLN
2	L	240	ASN
2	L	254	ASN

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Mol	Chain	Res	Type
2	L	276	ASN
2	L	380	ASN
2	L	383	ASN
2	L	454	ASN
1	M	34	GLN
1	M	94	GLN
1	M	98	GLN
2	N	240	ASN
2	N	254	ASN
2	N	270	GLN
2	N	277	ASN
2	N	361	GLN
2	N	380	ASN
2	N	383	ASN
2	N	428	ASN
2	N	437	ASN
2	N	454	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

12 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
3	NAG	N	800	2	14,14,15	0.50	0	17,19,21	0.78	0
3	NAG	H	800	2	14,14,15	0.47	0	17,19,21	1.23	1 (5%)
3	NAG	I	770	1	14,14,15	0.49	0	17,19,21	0.63	0
3	NAG	D	800	2	14,14,15	0.49	0	17,19,21	0.88	1 (5%)
3	NAG	F	800	2	14,14,15	0.43	0	17,19,21	1.24	1 (5%)
3	NAG	B	800	2	14,14,15	0.41	0	17,19,21	1.44	1 (5%)
3	NAG	E	770	1	14,14,15	0.37	0	17,19,21	1.78	3 (17%)
3	NAG	C	770	1	14,14,15	0.54	0	17,19,21	0.92	1 (5%)
3	NAG	G	770	1	14,14,15	0.60	0	17,19,21	1.33	1 (5%)
3	NAG	M	770	1	14,14,15	0.45	0	17,19,21	1.68	2 (11%)
3	NAG	A	770	1	14,14,15	0.47	0	17,19,21	0.77	1 (5%)
3	NAG	L	800	2	14,14,15	0.50	0	17,19,21	1.11	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
3	NAG	N	800	2	1/1/5/7	3/6/23/26	0/1/1/1
3	NAG	H	800	2	1/1/5/7	3/6/23/26	0/1/1/1
3	NAG	I	770	1	1/1/5/7	3/6/23/26	0/1/1/1
3	NAG	D	800	2	-	3/6/23/26	0/1/1/1
3	NAG	F	800	2	1/1/5/7	3/6/23/26	0/1/1/1
3	NAG	B	800	2	-	3/6/23/26	0/1/1/1
3	NAG	E	770	1	-	3/6/23/26	0/1/1/1
3	NAG	C	770	1	-	3/6/23/26	0/1/1/1
3	NAG	G	770	1	-	3/6/23/26	0/1/1/1
3	NAG	M	770	1	-	3/6/23/26	0/1/1/1
3	NAG	A	770	1	-	3/6/23/26	0/1/1/1
3	NAG	L	800	2	-	3/6/23/26	0/1/1/1

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	770	NAG	C1-O5-C5	5.75	119.99	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	M	770	NAG	C1-O5-C5	5.68	119.89	112.19
3	B	800	NAG	C1-O5-C5	4.78	118.67	112.19
3	G	770	NAG	C1-O5-C5	4.63	118.47	112.19
3	F	800	NAG	C1-O5-C5	4.33	118.06	112.19
3	H	800	NAG	C1-O5-C5	4.23	117.93	112.19
3	L	800	NAG	C1-O5-C5	3.91	117.49	112.19
3	M	770	NAG	O5-C1-C2	2.95	115.94	111.29
3	E	770	NAG	O5-C1-C2	2.37	115.04	111.29
3	D	800	NAG	C1-O5-C5	2.33	115.35	112.19
3	E	770	NAG	C4-C3-C2	-2.26	107.71	111.02
3	C	770	NAG	C1-O5-C5	2.09	115.03	112.19
3	A	770	NAG	C1-O5-C5	2.05	114.97	112.19

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	N	800	NAG	C1
3	H	800	NAG	C1
3	I	770	NAG	C1
3	F	800	NAG	C1

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	N	800	NAG	C1-C2-N2-C7
3	N	800	NAG	C8-C7-N2-C2
3	N	800	NAG	O7-C7-N2-C2
3	H	800	NAG	C8-C7-N2-C2
3	H	800	NAG	O7-C7-N2-C2
3	D	800	NAG	C8-C7-N2-C2
3	D	800	NAG	O7-C7-N2-C2
3	G	770	NAG	C8-C7-N2-C2
3	G	770	NAG	O7-C7-N2-C2
3	B	800	NAG	C8-C7-N2-C2
3	B	800	NAG	O7-C7-N2-C2
3	L	800	NAG	C8-C7-N2-C2
3	L	800	NAG	O7-C7-N2-C2
3	H	800	NAG	C1-C2-N2-C7
3	B	800	NAG	C1-C2-N2-C7
3	L	800	NAG	C1-C2-N2-C7
3	M	770	NAG	C8-C7-N2-C2
3	I	770	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	E	770	NAG	C8-C7-N2-C2
3	A	770	NAG	C8-C7-N2-C2
3	D	800	NAG	C1-C2-N2-C7
3	I	770	NAG	O7-C7-N2-C2
3	F	800	NAG	C8-C7-N2-C2
3	E	770	NAG	O7-C7-N2-C2
3	M	770	NAG	O7-C7-N2-C2
3	A	770	NAG	O7-C7-N2-C2
3	E	770	NAG	C1-C2-N2-C7
3	A	770	NAG	C1-C2-N2-C7
3	F	800	NAG	O7-C7-N2-C2
3	C	770	NAG	C8-C7-N2-C2
3	F	800	NAG	C3-C2-N2-C7
3	M	770	NAG	C3-C2-N2-C7
3	C	770	NAG	O7-C7-N2-C2
3	I	770	NAG	C3-C2-N2-C7
3	C	770	NAG	C3-C2-N2-C7
3	G	770	NAG	C3-C2-N2-C7

There are no ring outliers.

8 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	N	800	NAG	2	0
3	H	800	NAG	1	0
3	I	770	NAG	1	0
3	F	800	NAG	1	0
3	B	800	NAG	3	0
3	C	770	NAG	1	0
3	G	770	NAG	1	0
3	M	770	NAG	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	73/84 (86%)	-0.20	2 (2%) 54 44	25, 44, 85, 125	0
1	C	72/84 (85%)	-0.38	0 100 100	17, 35, 83, 157	0
1	E	73/84 (86%)	-0.21	1 (1%) 75 69	16, 40, 87, 121	0
1	G	72/84 (85%)	-0.11	1 (1%) 75 69	21, 50, 104, 169	0
1	I	71/84 (84%)	-0.30	0 100 100	12, 39, 88, 116	0
1	M	72/84 (85%)	-0.17	2 (2%) 53 43	15, 41, 86, 143	0
2	B	358/374 (95%)	-0.20	5 (1%) 75 69	14, 43, 82, 130	0
2	D	355/374 (94%)	-0.05	7 (1%) 65 56	15, 48, 95, 178	0
2	F	351/374 (93%)	-0.17	4 (1%) 80 75	14, 39, 87, 139	0
2	H	359/374 (95%)	-0.16	6 (1%) 70 63	15, 44, 84, 127	0
2	L	354/374 (94%)	-0.23	2 (0%) 89 86	13, 39, 82, 141	0
2	N	355/374 (94%)	-0.06	6 (1%) 70 63	17, 50, 94, 201	0
All	All	2565/2748 (93%)	-0.16	36 (1%) 75 69	12, 44, 91, 201	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	N	429	ARG	7.2
2	D	429	ARG	4.9
1	G	65	LYS	4.1
2	D	355	ALA	3.9
2	N	428	ASN	3.6
1	A	26	GLN	3.2
1	M	65	LYS	3.1
2	H	472	GLU	3.1
2	H	356	GLU	2.9
2	F	355	ALA	2.7
2	L	477	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
2	L	154	VAL	2.7
2	D	310	ASP	2.6
2	N	371	ASN	2.6
2	H	381	LEU	2.6
1	A	65	LYS	2.6
2	B	381	LEU	2.5
1	M	67	ASN	2.5
2	F	356	GLU	2.5
2	D	371	ASN	2.5
2	H	371	ASN	2.4
2	H	470	LYS	2.4
2	F	354	GLN	2.3
2	D	354	GLN	2.3
2	N	477	PHE	2.3
2	F	421	LYS	2.2
1	E	26	GLN	2.2
2	D	356	GLU	2.2
2	N	355	ALA	2.2
2	N	354	GLN	2.2
2	B	356	GLU	2.2
2	H	354	GLN	2.1
2	D	428	ASN	2.1
2	B	357	THR	2.1
2	B	419	LYS	2.1
2	B	470	LYS	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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. 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	B-factors(\AA^2)	Q<0.9
3	NAG	I	770	14/15	0.69	0.39	128,147,151,151	0
3	NAG	E	770	14/15	0.76	0.43	99,117,122,124	0
3	NAG	A	770	14/15	0.77	0.18	62,81,84,86	0
3	NAG	C	770	14/15	0.78	0.24	82,100,106,107	0
3	NAG	D	800	14/15	0.79	0.22	58,76,82,82	0
3	NAG	G	770	14/15	0.79	0.24	95,113,119,120	0
3	NAG	H	800	14/15	0.79	0.33	88,106,112,113	0
3	NAG	M	770	14/15	0.82	0.26	91,110,114,115	0
3	NAG	B	800	14/15	0.82	0.21	57,74,81,82	0
3	NAG	N	800	14/15	0.85	0.25	62,80,85,86	0
3	NAG	F	800	14/15	0.87	0.22	71,89,94,95	0
3	NAG	L	800	14/15	0.90	0.18	64,82,87,87	0

6.5 Other polymers [i](#)

There are no such residues in this entry.