



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:57 PM GMT

PDB ID : 3VI4
Title : Crystal structure of alpha5beta1 integrin headpiece in complex with RGD peptide
Authors : Nagae, M.; Nogi, T.; Takagi, J.
Deposited on : 2011-09-21
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

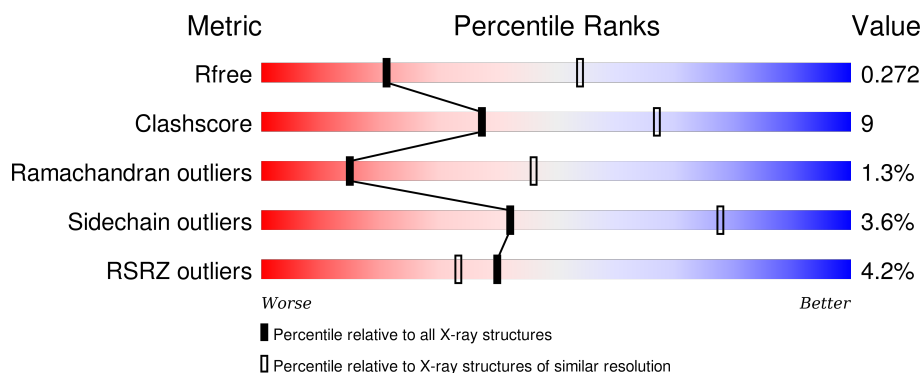
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	632	 3% 78% 16% • 5%
1	C	632	 5% 66% 26% • 7%
2	B	454	 9% 71% 22% • 5%
2	D	454	 6% 70% 24% • 5%
3	E	219	 % 74% 24% •

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Mol	Chain	Length	Quality of chain
3	L	219	<div><div></div><div>74%24%</div><div></div></div>
4	F	218	<div>%<div><div></div><div>74%24%</div><div></div></div></div>
4	H	218	<div>%<div><div></div><div>82%18%</div><div></div></div></div>
5	G	5	<div><div></div><div>60%20%20%</div><div></div></div>
5	I	5	<div><div></div><div>60%40%</div><div></div></div>

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 22882 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 Integrin alpha-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	601	Total	C	N	O	S	0	0	0
			4567	2892	768	893	14			
1	C	589	Total	C	N	O	S	0	0	0
			4466	2828	747	877	14			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	624	GLY	-	EXPRESSION TAG	UNP P08648
A	625	GLY	-	EXPRESSION TAG	UNP P08648
A	626	LEU	-	EXPRESSION TAG	UNP P08648
A	627	GLU	-	EXPRESSION TAG	UNP P08648
A	628	ASN	-	EXPRESSION TAG	UNP P08648
A	629	LEU	-	EXPRESSION TAG	UNP P08648
A	630	TYR	-	EXPRESSION TAG	UNP P08648
A	631	PHE	-	EXPRESSION TAG	UNP P08648
A	632	GLN	-	EXPRESSION TAG	UNP P08648
C	624	GLY	-	EXPRESSION TAG	UNP P08648
C	625	GLY	-	EXPRESSION TAG	UNP P08648
C	626	LEU	-	EXPRESSION TAG	UNP P08648
C	627	GLU	-	EXPRESSION TAG	UNP P08648
C	628	ASN	-	EXPRESSION TAG	UNP P08648
C	629	LEU	-	EXPRESSION TAG	UNP P08648
C	630	TYR	-	EXPRESSION TAG	UNP P08648
C	631	PHE	-	EXPRESSION TAG	UNP P08648
C	632	GLN	-	EXPRESSION TAG	UNP P08648

- Molecule 2 is a protein called Integrin beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	431	Total	C	N	O	S	0	0	0
			3353	2096	572	661	24			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	433	Total	C	N	O	S	0	0	0
			3372	2109	575	664	24			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	92	HIS	THR	SEE REMARK 999	UNP P05556
B	195	THR	SER	SEE REMARK 999	UNP P05556
B	446	GLY	-	EXPRESSION TAG	UNP P05556
B	447	GLY	-	EXPRESSION TAG	UNP P05556
B	448	LEU	-	EXPRESSION TAG	UNP P05556
B	449	GLU	-	EXPRESSION TAG	UNP P05556
B	450	ASN	-	EXPRESSION TAG	UNP P05556
B	451	LEU	-	EXPRESSION TAG	UNP P05556
B	452	TYR	-	EXPRESSION TAG	UNP P05556
B	453	PHE	-	EXPRESSION TAG	UNP P05556
B	454	GLN	-	EXPRESSION TAG	UNP P05556
D	92	HIS	THR	SEE REMARK 999	UNP P05556
D	195	THR	SER	SEE REMARK 999	UNP P05556
D	446	GLY	-	EXPRESSION TAG	UNP P05556
D	447	GLY	-	EXPRESSION TAG	UNP P05556
D	448	LEU	-	EXPRESSION TAG	UNP P05556
D	449	GLU	-	EXPRESSION TAG	UNP P05556
D	450	ASN	-	EXPRESSION TAG	UNP P05556
D	451	LEU	-	EXPRESSION TAG	UNP P05556
D	452	TYR	-	EXPRESSION TAG	UNP P05556
D	453	PHE	-	EXPRESSION TAG	UNP P05556
D	454	GLN	-	EXPRESSION TAG	UNP P05556

- Molecule 3 is a protein called SG/19 Fab fragment (Light chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	219	Total	C	N	O	S	0	0	0
			1701	1066	292	335	8			
3	E	219	Total	C	N	O	S	0	0	0
			1701	1066	292	335	8			

- Molecule 4 is a protein called SG/19 Fab fragment (Heavy chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	218	Total	C	N	O	S	0	0	0
			1651	1051	268	324	8			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	218	Total	C	N	O	S	0	0	0
			1651	1051	268	324	8			

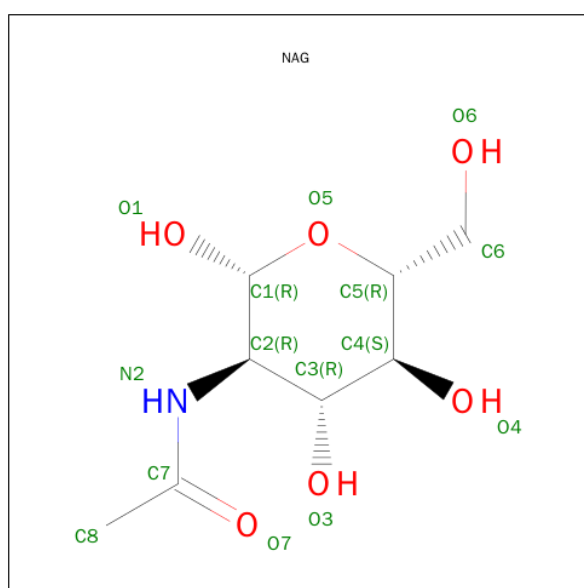
- Molecule 5 is a protein called RGD peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	G	4	Total	C	N	O	0	0	0
			31	16	8	7			
5	I	3	Total	C	N	O	0	0	0
			23	12	6	5			

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		
6	A	4	Total	Ca	0	0
			4	4		
6	D	1	Total	Ca	0	0
			1	1		
6	C	4	Total	Ca	0	0
			4	4		

- Molecule 7 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 8 is a polymer of unknown type called SUGAR (2-MER).

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

- Molecule 9 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	6	Total	C	N	O	0	0
			72	40	2	30		
9	C	6	Total	C	N	O	0	0
			72	40	2	30		

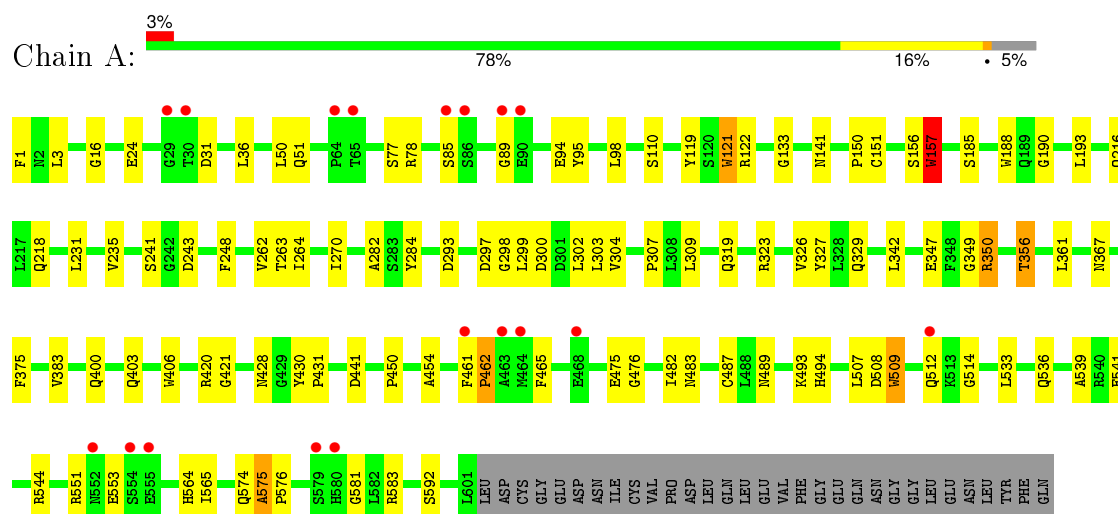
- Molecule 10 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	B	1	Total 1	Mg 1	0	0
10	D	1	Total 1	Mg 1	0	0

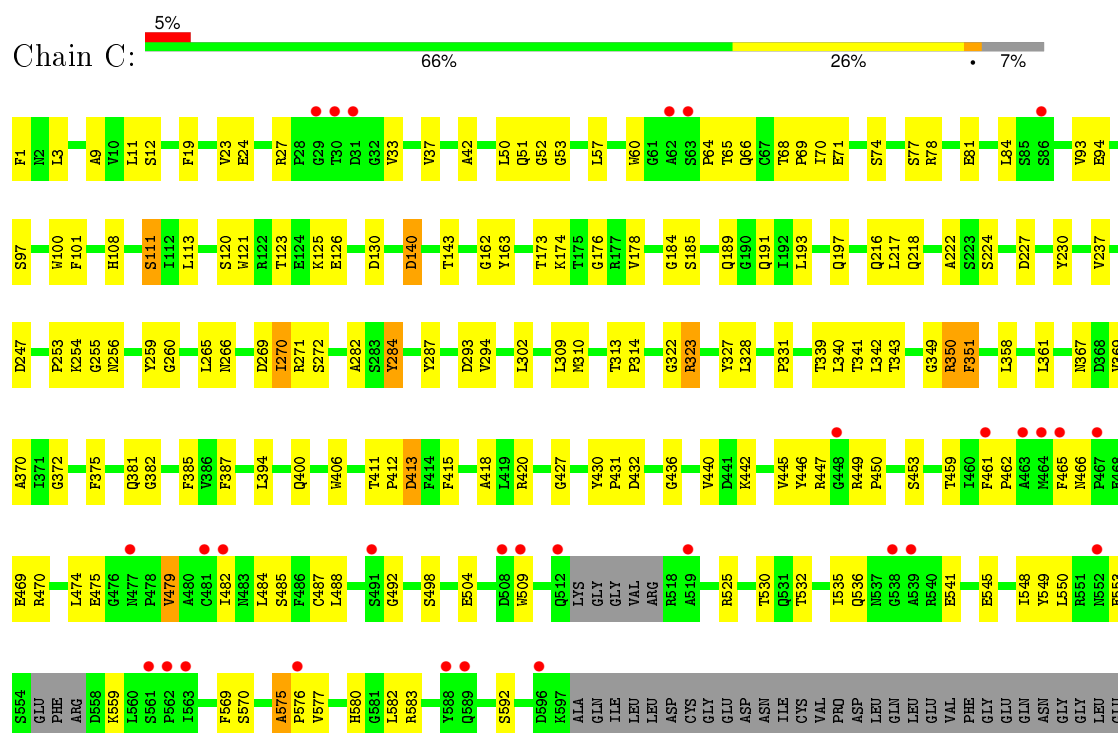
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Integrin alpha-5

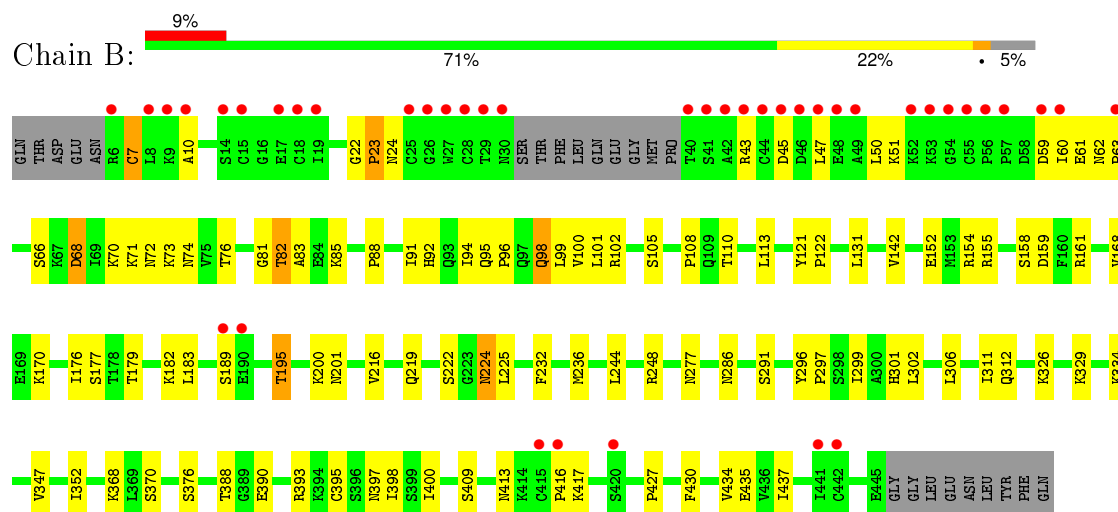


• Molecule 1: Integrin alpha-5

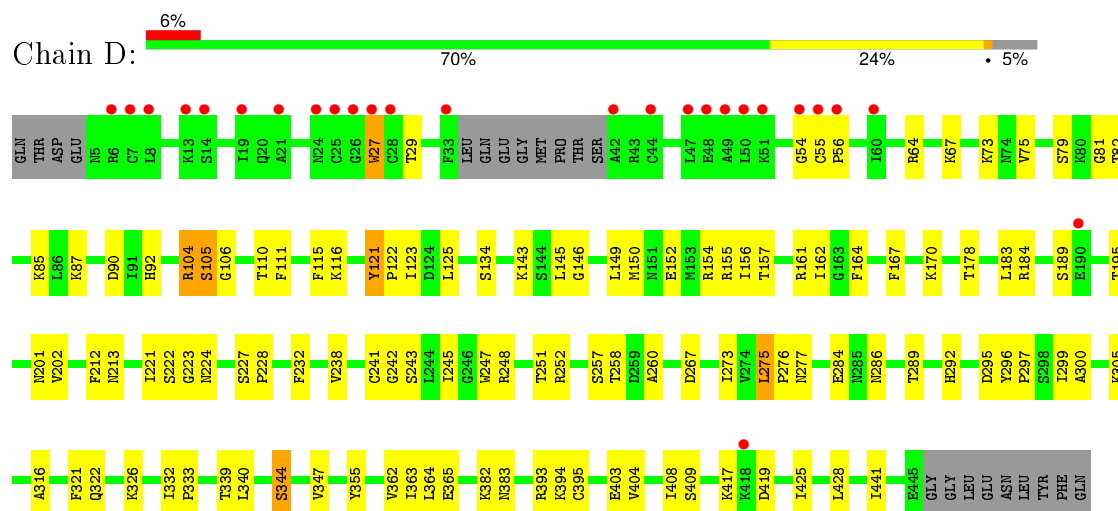


ASN
LEU
TYR
PHE
GLN

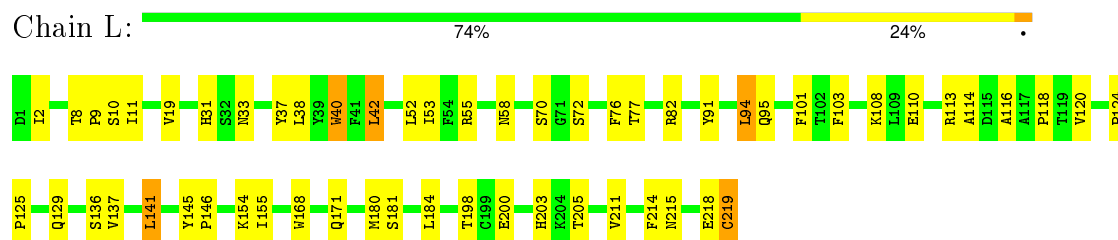
• Molecule 2: Integrin beta-1



• Molecule 2: Integrin beta-1

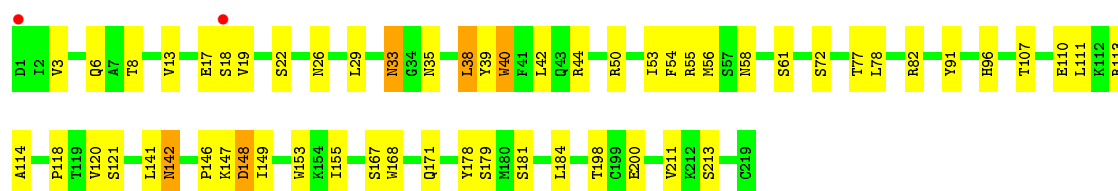


• Molecule 3: SG/19 Fab fragment (Light chain)

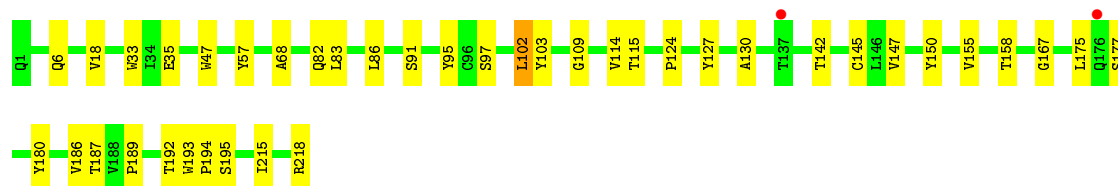
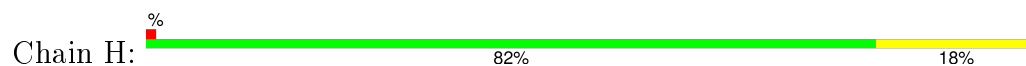


• Molecule 3: SG/19 Fab fragment (Light chain)

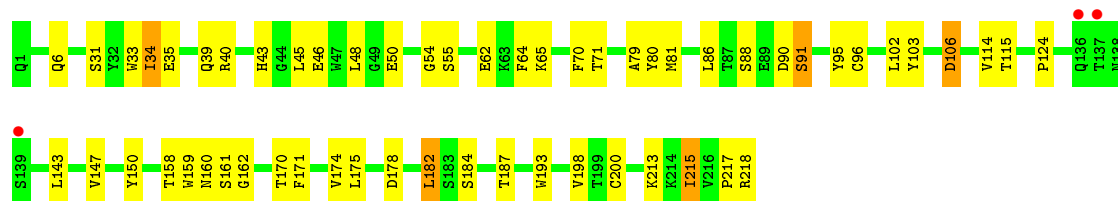
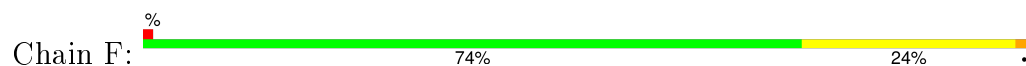




• Molecule 4: SG/19 Fab fragment (Heavy chain)



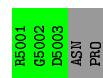
• Molecule 4: SG/19 Fab fragment (Heavy chain)



• Molecule 5: RGD peptide



• Molecule 5: RGD peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	92.81Å 103.87Å 125.26Å 75.72° 70.16° 70.90°	Depositor
Resolution (Å)	100.00 – 2.90 47.92 – 2.89	Depositor EDS
% Data completeness (in resolution range)	98.3 (100.00-2.90) 93.4 (47.92-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.51 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.215 , 0.275 0.213 , 0.272	Depositor DCC
R_{free} test set	4460 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	64.6	Xtriage
Anisotropy	0.474	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 31.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 89703 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	22882	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, MG, BMA, CA, 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	0.46	4/4684 (0.1%)	0.62	0/6374
1	C	0.48	4/4580 (0.1%)	0.63	0/6234
2	B	0.41	0/3410	0.59	0/4601
2	D	0.43	1/3430 (0.0%)	0.63	0/4628
3	E	0.50	3/1741 (0.2%)	0.65	1/2364 (0.0%)
3	L	0.51	2/1741 (0.1%)	0.63	1/2364 (0.0%)
4	F	0.56	1/1698 (0.1%)	0.63	0/2323
4	H	0.56	1/1698 (0.1%)	0.66	0/2323
5	G	0.49	0/30	0.52	0/38
5	I	0.49	0/22	0.57	0/27
All	All	0.48	16/23034 (0.1%)	0.62	2/31276 (0.0%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	40	TRP	CD2-CE2	5.51	1.48	1.41
1	C	406	TRP	CD2-CE2	5.48	1.48	1.41
3	E	40	TRP	CD2-CE2	5.46	1.47	1.41
1	A	121	TRP	CD2-CE2	5.43	1.47	1.41
3	L	168	TRP	CD2-CE2	5.34	1.47	1.41
1	A	188	TRP	CD2-CE2	5.20	1.47	1.41
3	E	153	TRP	CD2-CE2	5.19	1.47	1.41
1	A	157	TRP	CD2-CE2	5.17	1.47	1.41
4	F	193	TRP	CD2-CE2	5.17	1.47	1.41
1	C	100	TRP	CD2-CE2	5.12	1.47	1.41
1	C	509	TRP	CD2-CE2	5.11	1.47	1.41
1	A	406	TRP	CD2-CE2	5.08	1.47	1.41
2	D	27	TRP	CD2-CE2	5.07	1.47	1.41
4	H	47	TRP	CD2-CE2	5.06	1.47	1.41
1	C	60	TRP	CD2-CE2	5.02	1.47	1.41
3	E	168	TRP	CD2-CE2	5.01	1.47	1.41

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	94	LEU	CA-CB-CG	5.70	128.40	115.30
3	E	142	ASN	N-CA-C	5.38	125.53	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4567	0	4363	59	0
1	C	4466	0	4247	102	0
2	B	3353	0	3305	67	0
2	D	3372	0	3321	77	0
3	E	1701	0	1649	43	0
3	L	1701	0	1649	34	0
4	F	1651	0	1610	38	0
4	H	1651	0	1610	21	0
5	G	31	0	25	1	0
5	I	23	0	19	0	0
6	A	4	0	0	0	0
6	B	1	0	0	0	0
6	C	4	0	0	0	0
6	D	1	0	0	0	0
7	A	56	0	52	0	0
7	B	28	0	26	0	0
7	C	56	0	52	1	0
7	D	14	0	13	0	0
8	A	28	0	25	1	0
8	C	28	0	25	0	0
9	A	72	0	61	0	0
9	C	72	0	61	1	0
10	B	1	0	0	0	0
10	D	1	0	0	0	0
All	All	22882	0	22113	417	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 9.

All (417) 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:110:THR:HG22	2:B:409:SER:HB3	1.15	1.07
2:D:110:THR:HG22	2:D:409:SER:HB3	1.36	1.05
2:B:7:CYS:HB3	2:B:45:ASP:HA	1.39	1.01
2:D:284:GLU:HG2	2:D:289:THR:HG21	1.50	0.94
2:D:110:THR:HG22	2:D:409:SER:CB	1.98	0.93
2:B:110:THR:HG22	2:B:409:SER:CB	2.01	0.91
2:B:236:MET:HE3	2:B:301:HIS:HD2	1.36	0.89
2:B:152:GLU:HG2	2:B:352:ILE:HD11	1.54	0.88
1:C:189:GLN:HE22	1:C:222:ALA:H	1.19	0.88
2:B:68:ASP:HB3	2:B:70:LYS:HE2	1.55	0.86
3:E:147:LYS:O	3:E:147:LYS:HG3	1.76	0.86
1:A:475:GLU:HG3	1:A:476:GLY:H	1.40	0.85
2:B:71:LYS:HB3	2:B:98:GLN:HB2	1.59	0.84
2:B:158:SER:HA	3:E:55:ARG:HH12	1.43	0.83
3:E:54:PHE:HE2	3:E:55:ARG:HH21	1.27	0.82
1:A:241:SER:HB2	8:A:2008:NAG:H82	1.61	0.81
2:B:113:LEU:HD11	2:B:434:VAL:HG21	1.62	0.81
3:L:200:GLU:HG2	3:L:211:VAL:HG22	1.63	0.81
1:A:150:PRO:HB3	1:A:216:GLN:HE22	1.44	0.80
1:C:461:PHE:HB3	1:C:462:PRO:HD3	1.63	0.80
1:A:462:PRO:HD2	1:A:482:ILE:HG23	1.64	0.79
1:A:193:LEU:HD23	1:A:218:GLN:HG2	1.67	0.77
1:A:356:THR:HG21	1:A:421:GLY:H	1.47	0.77
1:C:504:GLU:HG3	1:C:530:THR:HG22	1.67	0.76
2:D:170:LYS:HD2	2:D:296:TYR:CZ	2.20	0.76
1:C:328:LEU:HD12	1:C:339:THR:HG21	1.68	0.75
2:D:161:ARG:NH2	2:D:248:ARG:HD3	2.00	0.75
3:E:33:ASN:HB3	3:E:35:ASN:HD22	1.52	0.75
2:B:45:ASP:H	2:B:50:LEU:HD12	1.52	0.75
1:C:78:ARG:HD3	1:C:94:GLU:OE2	1.87	0.74
2:B:152:GLU:CG	2:B:352:ILE:HD11	2.16	0.74
2:D:273:ILE:HD12	2:D:292:HIS:O	1.85	0.74
1:C:349:GLY:HA2	1:C:375:PHE:HB2	1.70	0.73
4:F:48:LEU:CD1	4:F:81:MET:HE1	2.18	0.73
2:B:236:MET:HE3	2:B:301:HIS:CD2	2.23	0.73
1:C:293:ASP:O	1:C:367:ASN:HB2	1.88	0.73
4:F:48:LEU:CD1	4:F:81:MET:CE	2.67	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:62:GLU:HA	4:F:65:LYS:HG3	1.71	0.73
1:A:509:TRP:O	1:A:512:GLN:HG2	1.89	0.72
1:C:323:ARG:NH1	1:C:343:THR:OG1	2.24	0.71
2:D:247:TRP:HB3	2:D:252:ARG:HE	1.56	0.71
2:B:236:MET:CE	2:B:301:HIS:HD2	2.03	0.71
1:A:293:ASP:O	1:A:367:ASN:HB2	1.90	0.71
2:D:162:ILE:HG21	2:D:212:PHE:CD1	2.26	0.71
2:D:247:TRP:HB3	2:D:252:ARG:NE	2.07	0.70
1:C:427:GLY:HA3	1:C:580:HIS:CE1	2.25	0.70
2:D:73:LYS:HD3	2:D:92:HIS:CE1	2.26	0.70
2:D:75:VAL:CG1	2:D:116:LYS:HD3	2.21	0.69
2:B:72:ASN:HB2	2:B:96:PRO:HA	1.74	0.68
1:C:53:GLY:HA3	1:C:101:PHE:O	1.92	0.68
4:F:48:LEU:HD13	4:F:81:MET:HE1	1.74	0.68
2:B:400:ILE:O	4:F:31:SER:HB3	1.95	0.67
1:A:24:GLU:HB3	1:A:36:LEU:HB2	1.77	0.67
3:E:142:ASN:O	3:E:179:SER:OG	2.10	0.67
1:A:77:SER:HB2	1:A:94:GLU:HG3	1.77	0.67
3:E:3:VAL:H	3:E:26:ASN:HB2	1.60	0.66
2:B:122:PRO:HB3	2:B:159:ASP:HB3	1.78	0.66
1:A:356:THR:HG21	1:A:421:GLY:N	2.09	0.66
1:C:569:PHE:HB2	1:C:592:SER:HB2	1.78	0.65
4:F:48:LEU:HD13	4:F:81:MET:CE	2.26	0.65
3:E:40:TRP:HB2	3:E:53:ILE:HB	1.78	0.65
3:E:13:VAL:HG21	3:E:19:VAL:HG21	1.78	0.64
2:D:201:ASN:HB3	2:D:286:ASN:HB3	1.79	0.64
1:A:575:ALA:N	1:A:576:PRO:HD3	2.12	0.64
1:C:453:SER:H	1:C:492:GLY:HA3	1.62	0.64
2:D:162:ILE:CG2	2:D:212:PHE:CD1	2.81	0.64
1:A:576:PRO:HG2	1:A:583:ARG:HG2	1.81	0.63
1:C:3:LEU:HB2	1:C:400:GLN:HE22	1.63	0.63
1:A:574:GLN:C	1:A:576:PRO:HD3	2.18	0.63
2:B:99:LEU:HD21	2:B:101:LEU:HD13	1.81	0.63
2:B:158:SER:HA	3:E:55:ARG:NH1	2.13	0.63
1:A:157:TRP:CZ2	2:B:225:LEU:HD11	2.34	0.63
2:D:104:ARG:HG3	2:D:105:SER:H	1.64	0.62
1:C:27:ARG:HG2	1:C:33:VAL:HG13	1.80	0.62
3:E:17:GLU:HG2	3:E:18:SER:H	1.64	0.62
1:C:256:ASN:HB2	1:C:259:TYR:HB2	1.82	0.62
2:D:284:GLU:CG	2:D:289:THR:HG21	2.27	0.62
1:C:351:PHE:O	1:C:372:GLY:O	2.17	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:PHE:H1	1:A:400:GLN:HB2	1.64	0.62
3:E:33:ASN:HB3	3:E:35:ASN:ND2	2.13	0.62
2:D:115:PHE:CE2	2:D:362:VAL:HG21	2.34	0.62
1:C:487:CYS:HB3	1:C:541:GLU:HB3	1.82	0.62
4:F:48:LEU:HD12	4:F:81:MET:HE3	1.82	0.61
4:F:46:GLU:OE2	4:F:64:PHE:HE1	1.83	0.61
2:B:236:MET:CE	2:B:301:HIS:CD2	2.81	0.61
2:D:85:LYS:HG2	3:L:37:TYR:CZ	2.35	0.60
2:B:176:ILE:HG22	2:B:224:ASN:HB2	1.83	0.60
2:D:156:ILE:HG21	2:D:355:TYR:CE2	2.36	0.60
3:L:120:VAL:HG13	3:L:141:LEU:HD23	1.83	0.60
3:E:8:THR:O	3:E:107:THR:HG22	2.01	0.60
4:F:48:LEU:HD12	4:F:81:MET:CE	2.29	0.60
3:E:146:PRO:HB2	3:E:148:ASP:HB2	1.84	0.60
2:D:178:THR:HA	2:D:183:LEU:HD13	1.82	0.60
3:E:13:VAL:HG21	3:E:19:VAL:CG2	2.30	0.60
4:F:70:PHE:CE1	4:F:81:MET:HG3	2.36	0.59
2:D:395:CYS:SG	2:D:404:VAL:HG11	2.42	0.59
3:L:113:ARG:HG2	3:L:114:ALA:N	2.17	0.59
4:H:142:THR:OG1	4:H:187:THR:HG22	2.01	0.59
2:B:201:ASN:HB3	2:B:286:ASN:HB3	1.83	0.59
1:C:575:ALA:HB3	1:C:576:PRO:HD3	1.83	0.59
1:C:411:THR:HB	1:C:412:PRO:HD2	1.84	0.59
1:A:110:SER:HB2	1:A:141:ASN:H	1.68	0.58
1:A:307:PRO:HB2	1:A:350:ARG:HB3	1.86	0.58
1:C:77:SER:HB2	1:C:94:GLU:HG3	1.86	0.58
1:C:113:LEU:HD11	1:C:178:VAL:HG13	1.84	0.58
2:B:142:VAL:HG12	2:B:216:VAL:HG11	1.86	0.58
2:B:121:TYR:CD2	2:B:122:PRO:HD2	2.39	0.57
3:E:6:GLN:NE2	3:E:107:THR:HG23	2.19	0.57
4:F:159:TRP:CZ3	4:F:200:CYS:HB3	2.39	0.57
4:F:143:LEU:HB3	4:F:215:ILE:HD12	1.86	0.57
1:C:111:SER:OG	1:C:197:GLN:NE2	2.37	0.57
2:B:110:THR:CG2	2:B:409:SER:HB3	2.10	0.57
1:C:485:SER:HB3	1:C:545:GLU:HG2	1.86	0.57
1:A:85:SER:HA	1:A:89:GLY:HA3	1.86	0.57
1:A:95:TYR:OH	1:A:122:ARG:HB2	2.05	0.57
4:F:91:SER:HA	4:F:114:VAL:O	2.05	0.57
1:C:361:LEU:HA	1:C:431:PRO:HG2	1.87	0.57
4:F:48:LEU:CD1	4:F:81:MET:HE3	2.34	0.56
1:A:487:CYS:HB3	1:A:541:GLU:HG2	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:323:ARG:HG2	1:C:342:LEU:O	2.05	0.56
1:C:382:GLY:HA3	1:C:415:PHE:O	2.05	0.56
2:B:83:ALA:HA	2:B:85:LYS:HZ3	1.70	0.56
3:L:118:PRO:HD3	3:L:203:HIS:ND1	2.21	0.56
1:C:140:ASP:HB2	1:C:143:THR:OG1	2.06	0.56
2:B:170:LYS:HE3	2:B:291:SER:O	2.06	0.56
2:D:164:PHE:HD2	2:D:202:VAL:HB	1.71	0.56
1:C:449:ARG:HG2	1:C:582:LEU:HB3	1.87	0.56
2:D:284:GLU:HG2	2:D:289:THR:CG2	2.31	0.56
2:D:152:GLU:HA	2:D:155:ARG:HH11	1.69	0.55
1:C:302:LEU:HB3	1:C:327:TYR:HB2	1.88	0.55
2:D:408:ILE:HD13	2:D:425:ILE:HD13	1.87	0.55
2:B:312:GLN:HG2	2:B:334:LYS:HG3	1.88	0.55
3:E:113:ARG:HG2	3:E:114:ALA:H	1.72	0.55
1:A:461:PHE:HB3	1:A:462:PRO:HD3	1.88	0.55
1:C:282:ALA:HB3	2:D:299:ILE:HD12	1.88	0.55
3:E:6:GLN:HE21	3:E:107:THR:HG23	1.73	0.54
4:F:39:GLN:HG3	4:F:45:LEU:HD23	1.88	0.54
3:E:111:LEU:H	3:E:171:GLN:HE22	1.55	0.54
1:C:78:ARG:O	1:C:93:VAL:HG12	2.07	0.54
3:E:17:GLU:HG2	3:E:18:SER:N	2.23	0.54
3:E:113:ARG:HG2	3:E:114:ALA:N	2.22	0.54
3:L:42:LEU:HD13	3:L:91:TYR:CZ	2.42	0.54
3:E:147:LYS:CG	3:E:147:LYS:O	2.51	0.54
1:C:272:SER:HB3	9:C:2010:NAG:H83	1.89	0.54
4:F:160:ASN:HD21	4:F:198:VAL:HA	1.71	0.54
2:D:110:THR:HG22	2:D:409:SER:HB2	1.87	0.54
2:B:82:THR:HG23	3:E:96:HIS:NE2	2.23	0.53
2:B:154:ARG:O	3:E:55:ARG:NH2	2.41	0.53
1:A:3:LEU:HB2	1:A:400:GLN:HE22	1.72	0.53
3:L:40:TRP:HB2	3:L:53:ILE:HB	1.91	0.53
2:B:236:MET:HE2	2:B:302:LEU:HD23	1.90	0.53
2:D:238:VAL:HA	2:D:245:ILE:HD13	1.90	0.53
2:B:82:THR:O	2:B:85:LYS:NZ	2.41	0.53
1:C:50:LEU:O	1:C:97:SER:HA	2.09	0.53
1:C:430:TYR:CE1	1:C:450:PRO:HA	2.43	0.53
1:A:193:LEU:CD2	1:A:218:GLN:HG2	2.37	0.53
1:C:313:THR:HB	1:C:314:PRO:HD2	1.89	0.53
2:D:154:ARG:NH1	3:L:58:ASN:OD1	2.41	0.53
3:E:54:PHE:O	3:E:55:ARG:HG3	2.09	0.53
1:A:536:GLN:HB3	1:A:539:ALA:HB2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:217:PRO:O	4:F:218:ARG:HB2	2.09	0.53
1:C:310:MET:HG3	2:D:300:ALA:HB2	1.89	0.53
4:F:62:GLU:HG2	4:F:65:LYS:HE3	1.92	0.52
2:D:67:LYS:NZ	2:D:111:PHE:HA	2.25	0.52
2:D:121:TYR:CD2	2:D:122:PRO:HD2	2.45	0.52
2:B:390:GLU:O	2:B:393:ARG:HG2	2.10	0.52
1:A:356:THR:HG21	1:A:420:ARG:HA	1.91	0.52
2:D:260:ALA:HA	2:D:321:PHE:CE2	2.44	0.52
1:C:12:SER:HB3	1:C:442:LYS:HG2	1.91	0.52
1:A:156:SER:O	1:A:157:TRP:HB2	2.10	0.51
2:D:104:ARG:HA	2:D:441:ILE:HB	1.92	0.51
1:A:302:LEU:HB3	1:A:327:TYR:HB2	1.92	0.51
1:C:81:GLU:O	1:C:84:LEU:HG	2.11	0.51
2:B:61:GLU:C	2:B:63:PRO:HD3	2.31	0.51
4:H:91:SER:HA	4:H:114:VAL:O	2.10	0.51
2:B:170:LYS:HD2	2:B:296:TYR:CZ	2.46	0.51
2:D:257:SER:HA	2:D:316:ALA:O	2.11	0.51
3:L:219:CYS:HB2	4:H:218:ARG:HH12	1.75	0.51
4:H:6:GLN:HE22	4:H:95:TYR:HA	1.75	0.51
1:C:331:PRO:O	3:E:82:ARG:NH1	2.44	0.50
1:C:474:LEU:HD23	1:C:475:GLU:N	2.26	0.50
1:C:349:GLY:O	1:C:351:PHE:N	2.44	0.50
4:F:40:ARG:HB2	4:F:43:HIS:HB2	1.93	0.50
3:E:120:VAL:HG22	3:E:141:LEU:HD22	1.92	0.50
1:A:509:TRP:CE3	1:A:564:HIS:HB3	2.46	0.50
2:B:74:ASN:HA	2:B:95:GLN:OE1	2.12	0.50
2:B:83:ALA:HA	2:B:85:LYS:NZ	2.26	0.49
1:A:297:ASP:O	1:A:299:LEU:N	2.34	0.49
1:C:488:LEU:HD12	1:C:535:ILE:HD11	1.94	0.49
1:C:71:GLU:OE1	1:C:74:SER:HB3	2.11	0.49
1:A:383:VAL:HG22	1:A:403:GLN:HG2	1.93	0.49
2:D:104:ARG:HG3	2:D:105:SER:N	2.27	0.49
3:L:55:ARG:NH1	4:H:102:LEU:O	2.45	0.49
2:B:105:SER:HB2	2:B:413:ASN:O	2.13	0.49
1:A:323:ARG:HG3	1:A:342:LEU:O	2.11	0.49
4:F:6:GLN:HE22	4:F:95:TYR:HA	1.77	0.49
2:B:98:GLN:HG2	2:B:435:GLU:HB2	1.93	0.49
1:C:465:PHE:CD1	1:C:550:LEU:HB2	2.47	0.49
1:C:254:LYS:HE3	2:D:267:ASP:OD2	2.13	0.49
1:C:570:SER:HB2	7:C:2016:NAG:O7	2.13	0.49
2:B:100:VAL:HG22	2:B:437:ILE:HB	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:269:ASP:O	1:C:270:ILE:C	2.51	0.49
2:B:47:LEU:HA	2:B:51:LYS:HD2	1.95	0.48
3:E:147:LYS:HB3	3:E:178:TYR:CZ	2.48	0.48
1:C:382:GLY:HA3	1:C:415:PHE:HB3	1.95	0.48
2:D:222:SER:OG	2:D:223:GLY:N	2.46	0.48
2:B:72:ASN:CB	2:B:96:PRO:HA	2.41	0.48
2:D:221:ILE:HG22	2:D:222:SER:O	2.13	0.48
1:C:163:TYR:O	1:C:184:GLY:HA3	2.13	0.48
1:C:418:ALA:HB3	1:C:436:GLY:HA3	1.95	0.48
3:E:53:ILE:HA	3:E:58:ASN:O	2.13	0.48
2:B:232:PHE:HB3	2:B:297:PRO:HD2	1.96	0.48
1:C:498:SER:HA	1:C:536:GLN:HG2	1.96	0.48
4:F:160:ASN:C	4:F:162:GLY:H	2.17	0.48
2:B:66:SER:H	2:B:102:ARG:HB2	1.78	0.48
1:A:493:LYS:HE3	1:A:494:HIS:CE1	2.49	0.48
3:E:155:ILE:HD11	3:E:184:LEU:HD21	1.95	0.48
3:E:42:LEU:HD13	3:E:91:TYR:CZ	2.49	0.48
1:A:303:LEU:HD22	1:A:326:VAL:HG22	1.95	0.48
1:C:475:GLU:OE2	1:C:525:ARG:NH2	2.46	0.48
2:B:326:LYS:O	2:B:329:LYS:HB3	2.14	0.48
1:C:173:THR:HG22	1:C:237:VAL:HG11	1.96	0.48
4:H:175:LEU:HD13	4:H:180:TYR:CE1	2.48	0.47
4:H:35:GLU:HB2	4:H:97:SER:HB3	1.96	0.47
2:D:184:ARG:HA	2:D:195:THR:HG22	1.95	0.47
1:C:19:PHE:CE1	1:C:37:VAL:HG11	2.49	0.47
1:C:224:SER:HB3	5:G:5001:ARG:HE	1.80	0.47
2:D:143:LYS:O	2:D:213:ASN:ND2	2.47	0.47
2:D:241:CYS:HB2	2:D:245:ILE:HD12	1.97	0.47
4:F:124:PRO:HB2	4:F:147:VAL:HG13	1.97	0.47
2:B:200:LYS:NZ	2:B:219:GLN:HE22	2.12	0.47
2:B:168:VAL:HG21	2:B:222:SER:HB3	1.97	0.47
2:D:75:VAL:HG11	2:D:116:LYS:HD3	1.95	0.47
1:C:385:PHE:HB3	1:C:387:PHE:CE1	2.49	0.47
2:D:322:GLN:O	2:D:326:LYS:HG2	2.15	0.47
1:A:507:LEU:HD23	1:A:565:ILE:HG23	1.96	0.47
1:C:265:LEU:HA	1:C:271:ARG:O	2.15	0.47
2:D:149:LEU:HD23	2:D:162:ILE:HD11	1.96	0.47
2:B:161:ARG:HH11	2:B:248:ARG:HE	1.62	0.47
1:C:255:GLY:H	1:C:260:GLY:HA2	1.79	0.46
3:L:95:GLN:NE2	3:L:101:PHE:HA	2.30	0.46
4:H:193:TRP:CD1	4:H:194:PRO:HA	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:129:GLN:HE22	3:L:136:SER:HB2	1.80	0.46
1:C:57:LEU:HB2	1:C:70:ILE:HD11	1.97	0.46
2:D:238:VAL:HA	2:D:245:ILE:CD1	2.45	0.46
2:B:179:THR:HB	2:B:182:LYS:HB2	1.97	0.46
1:A:190:GLY:HA3	1:A:231:LEU:HB3	1.96	0.46
1:C:474:LEU:HD23	1:C:475:GLU:H	1.79	0.46
4:F:170:THR:HG23	4:F:184:SER:HB2	1.96	0.46
2:D:116:LYS:HB2	2:D:403:GLU:HG2	1.98	0.46
1:C:19:PHE:CZ	1:C:37:VAL:HG11	2.50	0.46
3:E:44:ARG:NH1	3:E:50:ARG:HH11	2.13	0.46
1:A:282:ALA:HB3	2:B:299:ILE:HD12	1.96	0.46
3:L:113:ARG:HG2	3:L:114:ALA:O	2.16	0.46
1:C:465:PHE:CE1	1:C:550:LEU:HB2	2.51	0.46
1:C:140:ASP:HB2	1:C:143:THR:HG1	1.81	0.46
1:C:525:ARG:HG2	1:C:549:TYR:OH	2.16	0.46
4:F:124:PRO:HB3	4:F:150:TYR:HB3	1.98	0.46
4:H:18:VAL:HG12	4:H:86:LEU:HD11	1.98	0.46
2:D:167:PHE:CD1	2:D:167:PHE:C	2.89	0.46
2:B:50:LEU:HD22	2:B:51:LYS:HG3	1.98	0.46
2:B:73:LYS:HD3	2:B:92:HIS:CD2	2.51	0.46
2:D:152:GLU:HA	2:D:155:ARG:NH1	2.31	0.45
1:A:430:TYR:CE1	1:A:450:PRO:HA	2.51	0.45
1:A:551:ARG:C	1:A:553:GLU:H	2.17	0.45
3:L:31:HIS:HD2	3:L:33:ASN:H	1.63	0.45
2:D:156:ILE:HG21	2:D:355:TYR:HE2	1.81	0.45
2:B:427:PRO:HG2	2:B:430:PHE:HB2	1.97	0.45
1:C:577:VAL:HG22	1:C:583:ARG:HE	1.80	0.45
1:C:294:VAL:HG11	1:C:394:LEU:HB2	1.97	0.45
1:C:24:GLU:HA	1:C:420:ARG:HG2	1.98	0.45
1:C:130:ASP:HB3	1:C:162:GLY:O	2.17	0.45
3:E:198:THR:HG22	3:E:213:SER:OG	2.16	0.45
3:L:125:PRO:HD3	3:L:137:VAL:HG22	1.98	0.45
1:A:454:ALA:HA	1:A:489:ASN:O	2.16	0.45
1:C:340:LEU:HG	1:C:341:THR:N	2.30	0.45
3:E:110:GLU:OE2	3:E:147:LYS:HD2	2.16	0.45
2:B:62:ASN:N	2:B:63:PRO:HD3	2.32	0.45
3:L:154:LYS:HB2	3:L:198:THR:HB	1.98	0.45
2:D:162:ILE:CG2	2:D:212:PHE:HD1	2.29	0.45
3:E:22:SER:HB3	3:E:77:THR:HG22	1.99	0.44
3:L:110:GLU:HB3	3:L:171:GLN:HE22	1.82	0.44
2:D:340:LEU:HD11	2:D:344:SER:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:88:PRO:HA	2:B:91:ILE:HD12	1.99	0.44
4:F:35:GLU:OE2	4:F:50:GLU:OE1	2.35	0.44
4:F:86:LEU:HA	4:F:90:ASP:OD2	2.17	0.44
1:C:3:LEU:HD22	1:C:446:TYR:HB3	1.99	0.44
1:C:470:ARG:HD2	1:C:479:VAL:HA	1.99	0.44
2:D:395:CYS:SG	2:D:404:VAL:CG1	3.06	0.44
1:C:216:GLN:HG2	1:C:217:LEU:N	2.32	0.44
2:D:364:LEU:O	2:D:393:ARG:HD3	2.16	0.44
3:L:155:ILE:HD11	3:L:184:LEU:HD21	1.98	0.44
1:C:1:PHE:H1	1:C:400:GLN:HB2	1.83	0.44
4:F:91:SER:OG	4:F:115:THR:HA	2.18	0.44
1:A:428:ASN:HB2	1:A:430:TYR:HD1	1.82	0.44
4:H:130:ALA:HB2	4:H:215:ILE:HG23	2.00	0.44
4:F:71:THR:OG1	4:F:80:TYR:HB2	2.16	0.44
1:A:361:LEU:HA	1:A:431:PRO:HG2	2.00	0.44
1:A:509:TRP:HE3	1:A:564:HIS:HB3	1.83	0.44
3:L:124:PRO:HB3	3:L:214:PHE:CZ	2.53	0.44
3:E:38:LEU:HG	3:E:39:TYR:N	2.32	0.44
1:C:3:LEU:HD22	1:C:446:TYR:CB	2.48	0.44
2:D:305:LYS:HD3	2:D:305:LYS:HA	1.58	0.44
1:C:309:LEU:HD22	1:C:323:ARG:HB2	2.00	0.43
2:D:82:THR:O	2:D:85:LYS:NZ	2.48	0.43
1:C:413:ASP:HB3	1:C:440:VAL:HG22	2.00	0.43
2:D:125:LEU:C	2:D:125:LEU:HD23	2.38	0.43
2:D:75:VAL:HG11	2:D:116:LYS:HB3	2.00	0.43
1:A:309:LEU:HD22	1:A:323:ARG:HB2	2.00	0.43
4:H:167:GLY:O	4:H:186:VAL:HA	2.17	0.43
2:B:22:GLY:HA3	2:B:23:PRO:HD3	1.77	0.43
1:A:349:GLY:HA2	1:A:375:PHE:HB2	2.00	0.43
4:H:124:PRO:HB3	4:H:150:TYR:HB3	2.00	0.43
4:F:174:VAL:HG12	4:F:175:LEU:N	2.33	0.43
3:L:31:HIS:CD2	3:L:33:ASN:H	2.36	0.43
2:D:154:ARG:HB3	3:L:55:ARG:NH2	2.33	0.43
1:C:466:ASN:HB3	1:C:469:GLU:HB2	1.99	0.43
1:C:185:SER:HB2	1:C:191:GLN:HB2	2.01	0.43
2:B:131:LEU:HD23	2:B:131:LEU:HA	1.70	0.43
2:D:417:LYS:C	2:D:419:ASP:H	2.22	0.43
4:H:68:ALA:HA	4:H:82:GLN:O	2.19	0.43
1:C:322:GLY:HA3	1:C:351:PHE:HB3	2.00	0.43
2:D:363:ILE:HG12	2:D:394:LYS:HG2	2.00	0.43
1:A:248:PHE:O	1:A:264:ILE:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:475:GLU:HG3	1:A:476:GLY:N	2.19	0.43
3:E:3:VAL:H	3:E:26:ASN:CB	2.28	0.43
4:F:35:GLU:O	4:F:96:CYS:HA	2.19	0.43
3:L:94:LEU:HB2	3:L:103:PHE:CD1	2.54	0.43
2:D:85:LYS:HB2	3:L:31:HIS:CE1	2.53	0.43
1:C:465:PHE:CE1	1:C:482:ILE:HD13	2.54	0.43
1:C:432:ASP:OD1	1:C:447:ARG:HA	2.18	0.43
4:H:189:PRO:O	4:H:192:THR:OG1	2.35	0.43
1:C:23:VAL:HG23	1:C:420:ARG:HB2	2.01	0.42
1:C:358:LEU:HD11	1:C:370:ALA:HB2	2.00	0.42
4:H:102:LEU:HG	4:H:103:TYR:HB2	2.02	0.42
1:C:284:TYR:CD2	1:C:287:TYR:HD1	2.37	0.42
4:F:106:ASP:OD1	4:F:106:ASP:N	2.44	0.42
3:E:200:GLU:HG2	3:E:211:VAL:HG22	2.01	0.42
4:F:103:TYR:CE2	4:F:106:ASP:HB3	2.55	0.42
1:A:461:PHE:HB3	1:A:483:ASN:HB2	2.00	0.42
2:D:170:LYS:HD2	2:D:296:TYR:CE2	2.53	0.42
1:C:569:PHE:HD2	1:C:592:SER:HA	1.85	0.42
1:C:120:SER:HA	1:C:130:ASP:O	2.19	0.42
2:D:123:ILE:HD12	2:D:157:THR:HB	2.01	0.42
1:C:255:GLY:O	1:C:256:ASN:C	2.58	0.42
2:D:227:SER:HB3	2:D:228:PRO:HD3	2.01	0.42
1:A:50:LEU:HD23	1:A:98:LEU:HD21	2.01	0.42
2:B:236:MET:HE1	2:B:301:HIS:CD2	2.55	0.42
2:D:67:LYS:HZ2	2:D:111:PHE:HA	1.84	0.42
1:C:68:THR:HA	1:C:69:PRO:HD3	1.95	0.42
2:D:79:SER:HA	4:H:57:TYR:CD2	2.54	0.42
3:E:167:SER:OG	4:F:171:PHE:HB3	2.20	0.42
2:B:183:LEU:O	2:B:195:THR:HB	2.19	0.42
1:A:119:TYR:HB3	1:A:133:GLY:HA2	2.02	0.42
2:D:55:CYS:HA	2:D:56:PRO:HD3	1.83	0.42
3:L:52:LEU:C	3:L:53:ILE:HD12	2.40	0.41
2:B:155:ARG:NH1	3:E:61:SER:HB3	2.35	0.41
3:L:8:THR:HG21	3:L:11:ILE:HG12	2.02	0.41
2:D:365:GLU:HG2	2:D:428:LEU:HG	2.01	0.41
1:A:151:CYS:SG	1:A:193:LEU:HD11	2.60	0.41
1:A:50:LEU:O	1:A:51:GLN:C	2.59	0.41
3:L:180:MET:HG2	3:L:181:SER:N	2.36	0.41
1:A:319:GLN:HB2	1:A:347:GLU:HG2	2.01	0.41
1:A:300:ASP:O	1:A:329:GLN:NE2	2.49	0.41
4:H:124:PRO:HB2	4:H:147:VAL:HG13	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:VAL:HG21	1:A:304:VAL:CG2	2.50	0.41
1:A:24:GLU:OE2	1:A:24:GLU:HA	2.21	0.41
1:C:382:GLY:CA	1:C:415:PHE:HB3	2.50	0.41
3:E:113:ARG:NH1	3:E:114:ALA:O	2.51	0.41
2:D:81:GLY:HA2	4:H:33:TRP:CZ2	2.55	0.41
3:L:10:SER:HB2	3:L:108:LYS:HB3	2.02	0.41
2:D:382:LYS:HG2	2:D:383:ASN:HD22	1.86	0.41
1:C:53:GLY:H	1:C:101:PHE:H	1.68	0.41
2:B:142:VAL:CG1	2:B:142:VAL:O	2.67	0.41
2:D:232:PHE:HB3	2:D:297:PRO:HG2	2.02	0.41
1:C:230:TYR:HB2	1:C:253:PRO:HD2	2.02	0.41
2:B:395:CYS:HB3	2:B:398:ILE:HD11	2.03	0.41
3:L:215:ASN:HB2	3:L:218:GLU:HG3	2.03	0.41
2:D:275:LEU:HD12	2:D:276:PRO:HD2	2.01	0.41
2:D:277:ASN:HB2	2:D:295:ASP:O	2.21	0.41
1:C:484:LEU:HD12	1:C:548:ILE:HD13	2.03	0.41
2:D:87:LYS:HB2	2:D:90:ASP:OD1	2.21	0.41
1:C:42:ALA:O	1:C:52:GLY:O	2.39	0.41
1:C:427:GLY:HA3	1:C:580:HIS:HE1	1.81	0.41
3:L:116:ALA:O	3:L:205:THR:HG21	2.21	0.41
1:A:356:THR:CG2	1:A:421:GLY:H	2.26	0.41
4:H:6:GLN:HE21	4:H:109:GLY:HA3	1.86	0.41
4:F:150:TYR:OH	4:F:182:LEU:HD23	2.20	0.41
1:A:533:LEU:CD2	1:A:544:ARG:HE	2.33	0.41
2:D:332:ILE:HA	2:D:333:PRO:HD3	1.92	0.41
4:F:213:LYS:HD3	4:F:213:LYS:HA	1.90	0.41
1:C:125:LYS:HB2	1:C:126:GLU:H	1.53	0.41
2:B:81:GLY:HA2	4:F:33:TRP:CZ2	2.56	0.41
3:E:33:ASN:CB	3:E:35:ASN:HD22	2.28	0.41
3:L:145:TYR:CG	3:L:146:PRO:HA	2.56	0.41
2:B:7:CYS:HB3	2:B:45:ASP:CA	2.29	0.40
1:C:349:GLY:O	1:C:350:ARG:C	2.59	0.40
2:D:85:LYS:HG2	3:L:37:TYR:OH	2.21	0.40
1:A:16:GLY:N	1:A:441:ASP:OD1	2.50	0.40
1:C:247:ASP:OD1	1:C:266:ASN:HA	2.20	0.40
2:D:145:LEU:HG	2:D:146:GLY:N	2.35	0.40
4:H:175:LEU:HA	4:H:175:LEU:HD12	1.88	0.40
1:C:381:GLN:HB3	1:C:413:ASP:OD2	2.21	0.40
3:L:8:THR:HA	3:L:9:PRO:HD3	1.97	0.40
2:B:306:LEU:HD22	2:B:311:ILE:HB	2.03	0.40
1:C:108:HIS:CD2	1:C:176:GLY:CA	3.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:118:PRO:HG3	3:E:149:ILE:HD11	2.03	0.40
1:C:9:ALA:HB3	1:C:445:VAL:HB	2.03	0.40
4:F:215:ILE:HG12	4:F:215:ILE:H	1.78	0.40
3:L:129:GLN:HG3	4:H:127:TYR:CE2	2.57	0.40
3:E:29:LEU:HD12	3:E:38:LEU:HB2	2.03	0.40
4:F:34:ILE:HD13	4:F:79:ALA:HB2	2.04	0.40
1:C:193:LEU:HD23	1:C:218:GLN:HG2	2.04	0.40
3:L:38:LEU:HD22	3:L:76:PHE:CG	2.55	0.40
1:C:224:SER:HA	1:C:227:ASP:OD2	2.22	0.40
2:B:179:THR:HG22	2:B:182:LYS:H	1.87	0.40
2:B:417:LYS:HG2	2:B:417:LYS:H	1.72	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	599/632 (95%)	533 (89%)	56 (9%)	10 (2%)	11	38
1	C	583/632 (92%)	504 (86%)	69 (12%)	10 (2%)	11	38
2	B	427/454 (94%)	376 (88%)	46 (11%)	5 (1%)	16	48
2	D	429/454 (94%)	386 (90%)	38 (9%)	5 (1%)	16	48
3	E	217/219 (99%)	193 (89%)	22 (10%)	2 (1%)	21	57
3	L	217/219 (99%)	201 (93%)	13 (6%)	3 (1%)	14	44
4	F	216/218 (99%)	192 (89%)	22 (10%)	2 (1%)	21	57
4	H	216/218 (99%)	199 (92%)	16 (7%)	1 (0%)	34	71
5	G	2/5 (40%)	1 (50%)	1 (50%)	0	100	100
5	I	1/5 (20%)	1 (100%)	0	0	100	100
All	All	2907/3056 (95%)	2586 (89%)	283 (10%)	38 (1%)	15	46

All (38) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	575	ALA
2	B	416	PRO
1	C	350	ARG
1	C	575	ALA
1	A	514	GLY
1	A	581	GLY
2	B	10	ALA
2	B	43	ARG
2	B	189	SER
1	C	51	GLN
2	D	105	SER
4	F	54	GLY
1	A	78	ARG
1	A	350	ARG
2	B	23	PRO
1	C	553	GLU
1	C	559	LYS
2	D	54	GLY
2	D	189	SER
2	D	242	GLY
3	L	72	SER
3	E	56	MET
4	F	178	ASP
1	A	298	GLY
1	C	174	LYS
1	C	351	PHE
1	A	157	TRP
1	C	64	PRO
4	H	177	SER
3	E	33	ASN
1	A	243	ASP
1	A	462	PRO
1	C	140	ASP
2	D	106	GLY
3	L	82	ARG
1	A	270	ILE
1	C	270	ILE
3	L	2	ILE

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	486/513 (95%)	475 (98%)	11 (2%)	58	87
1	C	476/513 (93%)	463 (97%)	13 (3%)	52	84
2	B	382/402 (95%)	361 (94%)	21 (6%)	27	61
2	D	384/402 (96%)	369 (96%)	15 (4%)	39	75
3	E	194/194 (100%)	188 (97%)	6 (3%)	47	82
3	L	194/194 (100%)	188 (97%)	6 (3%)	47	82
4	F	186/186 (100%)	175 (94%)	11 (6%)	24	58
4	H	186/186 (100%)	179 (96%)	7 (4%)	40	76
5	G	3/4 (75%)	3 (100%)	0	100	100
5	I	2/4 (50%)	2 (100%)	0	100	100
All	All	2493/2598 (96%)	2403 (96%)	90 (4%)	42	78

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ASP
1	A	121	TRP
1	A	185	SER
1	A	235	VAL
1	A	263	THR
1	A	284	TYR
1	A	356	THR
1	A	465	PHE
1	A	508	ASP
1	A	509	TRP
1	A	592	SER
2	B	7	CYS
2	B	24	ASN
2	B	59	ASP
2	B	60	ILE
2	B	68	ASP

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Mol	Chain	Res	Type
2	B	76	THR
2	B	82	THR
2	B	94	ILE
2	B	98	GLN
2	B	108	PRO
2	B	177	SER
2	B	195	THR
2	B	224	ASN
2	B	244	LEU
2	B	277	ASN
2	B	347	VAL
2	B	368	LYS
2	B	370	SER
2	B	376	SER
2	B	388	THR
2	B	397	ASN
1	C	11	LEU
1	C	65	THR
1	C	66	GLN
1	C	111	SER
1	C	121	TRP
1	C	123	THR
1	C	284	TYR
1	C	323	ARG
1	C	369	VAL
1	C	413	ASP
1	C	459	THR
1	C	479	VAL
1	C	532	THR
2	D	27	TRP
2	D	29	THR
2	D	64	ARG
2	D	104	ARG
2	D	121	TYR
2	D	134	SER
2	D	150	MET
2	D	224	ASN
2	D	243	SER
2	D	251	THR
2	D	258	THR
2	D	275	LEU
2	D	339	THR

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Mol	Chain	Res	Type
2	D	344	SER
2	D	347	VAL
3	L	19	VAL
3	L	42	LEU
3	L	70	SER
3	L	77	THR
3	L	141	LEU
3	L	219	CYS
4	H	83	LEU
4	H	102	LEU
4	H	115	THR
4	H	145	CYS
4	H	155	VAL
4	H	158	THR
4	H	195	SER
3	E	38	LEU
3	E	72	SER
3	E	78	LEU
3	E	121	SER
3	E	148	ASP
3	E	181	SER
4	F	34	ILE
4	F	55	SER
4	F	88	SER
4	F	91	SER
4	F	102	LEU
4	F	106	ASP
4	F	158	THR
4	F	161	SER
4	F	182	LEU
4	F	187	THR
4	F	215	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	108	HIS
1	A	161	GLN
1	A	165	GLN
1	A	214	GLN
1	A	216	GLN

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Mol	Chain	Res	Type
1	A	400	GLN
1	A	477	ASN
1	A	494	HIS
2	B	92	HIS
2	B	151	ASN
2	B	219	GLN
2	B	277	ASN
2	B	292	HIS
2	B	301	HIS
2	B	310	ASN
2	B	405	GLN
1	C	108	HIS
1	C	165	GLN
1	C	189	GLN
1	C	197	GLN
1	C	400	GLN
1	C	552	ASN
1	C	574	GLN
1	C	580	HIS
2	D	5	ASN
2	D	92	HIS
2	D	95	GLN
2	D	213	ASN
2	D	312	GLN
2	D	383	ASN
2	D	405	GLN
3	L	31	HIS
3	L	43	GLN
3	L	171	GLN
4	H	6	GLN
4	H	39	GLN
3	E	26	ASN
3	E	33	ASN
3	E	35	ASN
3	E	171	GLN
3	E	194	HIS
4	F	6	GLN
4	F	39	GLN
4	F	160	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

16 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$
8	NAG	A	2008	1,8	14,14,15	0.47	0	15,19,21	2.14	4 (26%)
8	NAG	A	2009	8	14,14,15	0.49	0	15,19,21	1.43	2 (13%)
9	NAG	A	2010	1,9	14,14,15	0.71	0	15,19,21	0.84	0
9	NAG	A	2011	9	14,14,15	0.59	0	15,19,21	2.59	4 (26%)
9	BMA	A	2012	9	11,11,12	0.50	0	14,15,17	0.95	0
9	MAN	A	2013	9	11,11,12	0.66	0	14,15,17	0.71	0
9	MAN	A	2014	9	11,11,12	0.57	0	14,15,17	1.36	2 (14%)
9	MAN	A	2015	9	11,11,12	0.81	0	14,15,17	1.59	3 (21%)
8	NAG	C	2005	1,8	14,14,15	0.49	0	15,19,21	2.49	4 (26%)
8	NAG	C	2006	8	14,14,15	0.49	0	15,19,21	0.78	0
9	NAG	C	2010	9,1	14,14,15	0.67	0	15,19,21	1.25	2 (13%)
9	NAG	C	2011	9	14,14,15	0.54	0	15,19,21	2.77	3 (20%)
9	BMA	C	2012	9	11,11,12	0.40	0	14,15,17	1.03	1 (7%)
9	MAN	C	2013	9	11,11,12	0.64	0	14,15,17	1.22	1 (7%)
9	MAN	C	2014	9	11,11,12	0.60	0	14,15,17	1.62	4 (28%)
9	MAN	C	2015	9	11,11,12	0.52	0	14,15,17	1.01	1 (7%)

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
8	NAG	A	2008	1,8	-	0/6/23/26	0/1/1/1
8	NAG	A	2009	8	-	0/6/23/26	0/1/1/1
9	NAG	A	2010	1,9	-	0/6/23/26	0/1/1/1
9	NAG	A	2011	9	-	0/6/23/26	0/1/1/1
9	BMA	A	2012	9	-	0/2/19/22	0/1/1/1
9	MAN	A	2013	9	-	0/2/19/22	0/1/1/1
9	MAN	A	2014	9	-	0/2/19/22	0/1/1/1
9	MAN	A	2015	9	-	0/2/19/22	0/1/1/1
8	NAG	C	2005	1,8	-	0/6/23/26	0/1/1/1
8	NAG	C	2006	8	-	0/6/23/26	0/1/1/1
9	NAG	C	2010	9,1	-	0/6/23/26	0/1/1/1
9	NAG	C	2011	9	-	0/6/23/26	0/1/1/1
9	BMA	C	2012	9	-	0/2/19/22	0/1/1/1
9	MAN	C	2013	9	-	0/2/19/22	0/1/1/1
9	MAN	C	2014	9	-	0/2/19/22	0/1/1/1
9	MAN	C	2015	9	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	2011	NAG	C4-C3-C2	-5.16	103.21	111.23
9	C	2011	NAG	C4-C3-C2	-4.66	103.98	111.23
8	C	2005	NAG	C4-C3-C2	-4.53	104.18	111.23
8	A	2009	NAG	C2-N2-C7	-2.90	119.32	123.04
8	A	2008	NAG	C6-C5-C4	-2.48	106.90	113.02
9	C	2012	BMA	C1-O5-C5	-2.39	109.22	112.25
9	C	2010	NAG	O4-C4-C3	-2.32	105.11	110.34
9	C	2014	MAN	O5-C5-C6	2.06	111.80	107.35
9	C	2014	MAN	C1-C2-C3	2.11	112.03	109.54
9	A	2011	NAG	O3-C3-C2	2.12	113.32	109.11
8	A	2008	NAG	C3-C4-C5	2.20	114.03	110.20
9	A	2014	MAN	O5-C5-C6	2.28	112.29	107.35
9	C	2010	NAG	C2-N2-C7	2.62	126.40	123.04
9	C	2014	MAN	C2-C3-C4	2.74	115.70	111.04
9	C	2013	MAN	C1-C2-C3	2.90	112.97	109.54
9	C	2014	MAN	C3-C4-C5	2.92	115.29	110.20
8	C	2005	NAG	C3-C2-N2	2.93	117.59	110.56
9	C	2015	MAN	C1-O5-C5	2.97	116.02	112.25
9	A	2015	MAN	C1-C2-C3	3.07	113.17	109.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	2005	NAG	C2-N2-C7	3.17	127.11	123.04
9	A	2015	MAN	C3-C4-C5	3.33	116.00	110.20
9	A	2015	MAN	C2-C3-C4	3.41	116.83	111.04
8	A	2008	NAG	C4-C3-C2	3.50	116.67	111.23
8	A	2009	NAG	C1-O5-C5	3.68	116.92	112.25
9	A	2014	MAN	C1-O5-C5	3.82	117.09	112.25
9	C	2011	NAG	C3-C2-N2	5.04	122.62	110.56
9	A	2011	NAG	C3-C2-N2	5.26	123.17	110.56
8	A	2008	NAG	C1-O5-C5	5.46	119.18	112.25
9	A	2011	NAG	C2-N2-C7	6.02	130.77	123.04
8	C	2005	NAG	C1-O5-C5	6.43	120.41	112.25
9	C	2011	NAG	C2-N2-C7	7.77	133.02	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	2008	NAG	1	0
9	C	2010	NAG	1	0

5.6 Ligand geometry ⓘ

Of 23 ligands modelled in this entry, 12 are monoatomic - leaving 11 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
7	NAG	A	2005	1	14,14,15	0.47	0	15,19,21	1.17	1 (6%)
7	NAG	A	2006	1	14,14,15	0.56	0	15,19,21	1.10	1 (6%)
7	NAG	A	2007	1	14,14,15	0.48	0	15,19,21	0.82	0
7	NAG	A	2016	1	14,14,15	0.48	0	15,19,21	2.25	2 (13%)
7	NAG	B	503	2	14,14,15	0.55	0	15,19,21	1.00	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	NAG	B	504	2	14,14,15	0.45	0	15,19,21	1.09	1 (6%)
7	NAG	C	2007	1	14,14,15	0.51	0	15,19,21	0.90	1 (6%)
7	NAG	C	2008	1	14,14,15	0.51	0	15,19,21	1.03	1 (6%)
7	NAG	C	2009	1	14,14,15	0.51	0	15,19,21	1.32	1 (6%)
7	NAG	C	2016	1	14,14,15	0.43	0	15,19,21	1.41	1 (6%)
7	NAG	D	503	2	14,14,15	0.36	0	15,19,21	2.03	2 (13%)

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
7	NAG	A	2005	1	-	0/6/23/26	0/1/1/1
7	NAG	A	2006	1	-	0/6/23/26	0/1/1/1
7	NAG	A	2007	1	-	0/6/23/26	0/1/1/1
7	NAG	A	2016	1	-	0/6/23/26	0/1/1/1
7	NAG	B	503	2	-	0/6/23/26	0/1/1/1
7	NAG	B	504	2	-	0/6/23/26	0/1/1/1
7	NAG	C	2007	1	-	0/6/23/26	0/1/1/1
7	NAG	C	2008	1	-	0/6/23/26	0/1/1/1
7	NAG	C	2009	1	-	0/6/23/26	0/1/1/1
7	NAG	C	2016	1	-	0/6/23/26	0/1/1/1
7	NAG	D	503	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	503	NAG	C6-C5-C4	-2.41	107.07	113.02
7	A	2016	NAG	C3-C4-C5	-2.04	106.64	110.20
7	A	2006	NAG	C2-N2-C7	2.25	125.92	123.04
7	C	2007	NAG	C1-O5-C5	2.36	115.24	112.25
7	B	503	NAG	C1-O5-C5	2.47	115.38	112.25
7	C	2008	NAG	C1-O5-C5	2.89	115.92	112.25
7	B	504	NAG	C1-O5-C5	3.11	116.19	112.25
7	C	2009	NAG	C1-O5-C5	3.67	116.90	112.25
7	A	2005	NAG	C1-O5-C5	3.73	116.98	112.25
7	C	2016	NAG	C1-O5-C5	4.45	117.90	112.25
7	D	503	NAG	C1-O5-C5	6.66	120.69	112.25
7	A	2016	NAG	C1-O5-C5	7.63	121.93	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	2016	NAG	1	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	601/632 (95%)	-0.04	18 (2%) 54 47	40, 66, 115, 159	0
1	C	589/632 (93%)	0.10	30 (5%) 32 25	37, 65, 131, 164	0
2	B	431/454 (94%)	0.31	41 (9%) 10 6	41, 64, 161, 217	0
2	D	433/454 (95%)	-0.02	26 (6%) 25 18	36, 58, 139, 160	0
3	E	219/219 (100%)	-0.35	2 (0%) 85 84	42, 62, 80, 94	0
3	L	219/219 (100%)	-0.43	0 100 100	38, 55, 73, 85	0
4	F	218/218 (100%)	-0.22	3 (1%) 78 76	44, 66, 106, 127	0
4	H	218/218 (100%)	-0.33	2 (0%) 85 84	35, 58, 97, 111	0
5	G	4/5 (80%)	0.04	0 100 100	85, 87, 88, 99	0
5	I	3/5 (60%)	-0.23	0 100 100	77, 77, 82, 86	0
All	All	2935/3056 (96%)	-0.05	122 (4%) 40 33	35, 62, 127, 217	0

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	43	ARG	13.6
2	B	42	ALA	13.5
2	B	41	SER	12.3
2	B	28	CYS	9.7
2	B	19	ILE	7.9
1	A	64	PRO	7.2
1	C	464	MET	7.0
2	D	33	PHE	5.9
2	B	18	CYS	5.7
2	D	42	ALA	5.6
2	B	27	TRP	5.5
2	B	40	THR	5.5
1	C	477	ASN	5.5

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Mol	Chain	Res	Type	RSRZ
1	C	465	PHE	5.3
1	A	512	GLN	5.3
4	F	137	THR	5.2
1	A	579	SER	5.2
1	A	580	HIS	5.0
1	A	30	THR	5.0
2	B	47	LEU	4.8
2	B	45	ASP	4.8
1	C	561	SER	4.7
2	B	49	ALA	4.7
2	B	30	ASN	4.7
2	D	49	ALA	4.5
2	B	60	ILE	4.5
2	B	8	LEU	4.5
1	C	519	ALA	4.5
2	B	15	CYS	4.4
2	B	44	CYS	4.4
2	D	54	GLY	4.3
2	D	51	LYS	4.3
2	D	50	LEU	4.3
2	B	25	CYS	4.2
1	C	62	ALA	4.2
1	C	29	GLY	4.2
2	B	56	PRO	4.2
2	D	190	GLU	4.2
2	B	416	PRO	4.1
2	D	27	TRP	4.1
2	B	10	ALA	4.0
2	B	17	GLU	3.9
2	D	7	CYS	3.9
2	B	441	ILE	3.9
1	C	588	TYR	3.9
1	A	86	SER	3.8
1	C	576	PRO	3.8
1	C	539	ALA	3.7
2	D	47	LEU	3.7
2	D	26	GLY	3.7
2	D	418	LYS	3.7
2	B	14	SER	3.5
1	A	555	GLU	3.5
2	D	24	ASN	3.5
1	A	89	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
2	B	54	GLY	3.4
2	B	52	LYS	3.3
2	B	53	LYS	3.3
2	B	190	GLU	3.3
2	B	442	CYS	3.3
1	C	63	SER	3.3
1	C	589	GLN	3.1
2	B	55	CYS	3.1
2	D	13	LYS	3.0
2	D	44	CYS	3.0
1	A	90	GLU	3.0
1	C	482	ILE	2.9
2	D	6	ARG	2.9
1	C	562	PRO	2.9
1	C	461	PHE	2.9
2	B	9	LYS	2.8
1	C	512	GLN	2.8
2	B	46	ASP	2.8
1	A	29	GLY	2.8
1	C	509	TRP	2.7
1	A	554	SER	2.7
1	C	481	CYS	2.7
2	B	57	PRO	2.7
1	A	468	GLU	2.6
2	B	48	GLU	2.6
2	D	8	LEU	2.6
2	B	59	ASP	2.6
1	C	86	SER	2.6
2	D	14	SER	2.6
1	A	464	MET	2.6
3	E	1	ASP	2.6
2	D	25	CYS	2.5
2	B	63	PRO	2.5
2	D	60	ILE	2.5
4	H	176	GLN	2.5
1	A	552	ASN	2.5
1	C	467	PRO	2.5
1	A	463	ALA	2.5
1	C	552	ASN	2.4
1	C	596	ASP	2.4
1	A	65	THR	2.4
2	D	19	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	26	GLY	2.4
2	B	189	SER	2.4
2	B	415	CYS	2.4
1	C	491	SER	2.4
1	C	463	ALA	2.3
2	B	420	SER	2.3
2	D	28	CYS	2.3
1	A	461	PHE	2.3
2	B	29	THR	2.3
2	D	56	PRO	2.3
4	F	139	SER	2.2
1	C	538	GLY	2.2
4	F	136	GLN	2.2
1	C	563	ILE	2.1
2	D	21	ALA	2.1
1	C	30	THR	2.1
2	D	48	GLU	2.1
4	H	137	THR	2.1
1	C	508	ASP	2.1
1	C	31	ASP	2.1
2	D	55	CYS	2.1
1	A	85	SER	2.1
2	B	6	ARG	2.0
1	C	448	GLY	2.0
3	E	18	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(Å ²)	Q<0.9
8	NAG	A	2008	14/15	0.93	0.23	1.45	74,81,86,92	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
8	NAG	C	2005	14/15	0.91	0.19	0.87	73,78,87,88	0
9	NAG	C	2010	14/15	0.96	0.11	-2.10	50,54,60,64	0
9	NAG	A	2010	14/15	0.96	0.11	-2.10	54,59,62,68	0
8	NAG	C	2006	14/15	0.88	0.24	-	88,94,98,100	0
9	BMA	C	2012	11/12	0.92	0.10	-	71,76,82,85	0
9	MAN	A	2014	11/12	0.86	0.18	-	85,95,103,104	0
9	MAN	A	2015	11/12	0.90	0.11	-	84,89,94,97	0
9	MAN	C	2015	11/12	0.88	0.15	-	79,85,97,103	0
9	NAG	C	2011	14/15	0.94	0.16	-	57,61,68,77	0
9	BMA	A	2012	11/12	0.84	0.13	-	72,80,84,86	0
9	NAG	A	2011	14/15	0.93	0.16	-	61,64,71,75	0
8	NAG	A	2009	14/15	0.79	0.31	-	85,99,103,104	0
9	MAN	A	2013	11/12	0.90	0.14	-	78,84,87,95	0
9	MAN	C	2013	11/12	0.78	0.16	-	81,91,95,96	0
9	MAN	C	2014	11/12	0.82	0.12	-	83,86,92,97	0

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. 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
10	MG	B	502	1/1	0.94	0.21	0.78	66,66,66,66	0
7	NAG	A	2005	14/15	0.90	0.17	0.19	64,72,79,79	0
6	CA	A	2001	1/1	0.95	0.13	0.09	69,69,69,69	0
7	NAG	A	2016	14/15	0.85	0.19	0.05	84,89,95,99	0
7	NAG	C	2016	14/15	0.79	0.19	-0.26	88,94,106,111	0
6	CA	A	2003	1/1	0.93	0.13	-0.45	73,73,73,73	0
6	CA	C	2001	1/1	0.95	0.12	-0.73	58,58,58,58	0
6	CA	C	2004	1/1	0.87	0.11	-0.79	76,76,76,76	0
7	NAG	B	503	14/15	0.86	0.15	-0.82	60,62,67,69	0
6	CA	A	2004	1/1	0.79	0.13	-0.96	81,81,81,81	0
6	CA	C	2003	1/1	0.95	0.10	-1.39	78,78,78,78	0
6	CA	D	501	1/1	0.93	0.12	-1.83	50,50,50,50	0
10	MG	D	502	1/1	0.80	0.10	-2.08	66,66,66,66	0
6	CA	A	2002	1/1	0.92	0.06	-2.14	81,81,81,81	0
6	CA	C	2002	1/1	0.97	0.07	-3.24	80,80,80,80	0
6	CA	B	501	1/1	0.99	0.09	-3.81	59,59,59,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
7	NAG	A	2007	14/15	0.90	0.14	-	78,88,93,97	0
7	NAG	B	504	14/15	0.90	0.15	-	67,70,76,81	0
7	NAG	C	2007	14/15	0.74	0.18	-	67,78,85,86	0
7	NAG	A	2006	14/15	0.77	0.16	-	98,103,109,113	0
7	NAG	C	2008	14/15	0.77	0.18	-	103,112,114,114	0
7	NAG	D	503	14/15	0.83	0.14	-	55,62,65,66	0
7	NAG	C	2009	14/15	0.92	0.14	-	60,67,71,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.