



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 8, 2018 – 05:14 PM EST

PDB ID : 5VQF
Title : Crystal Structure of pro-TGF-beta 1
Authors : Zhao, B.; Xu, S.; Dong, X.; Lu, C.; Springer, T.A.
Deposited on : 2017-05-08
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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030736

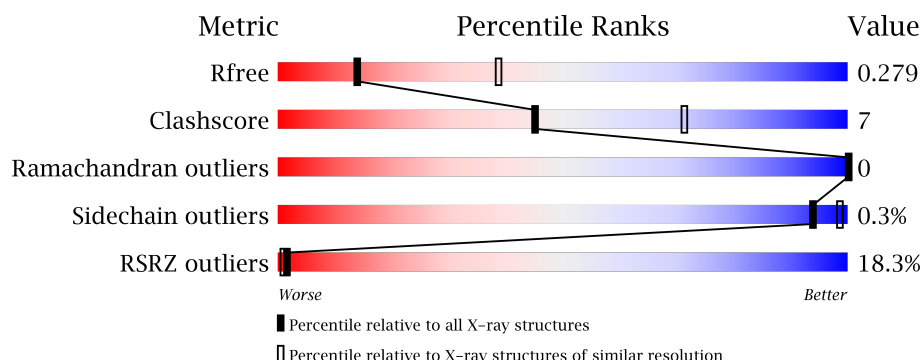
1 Overall quality at a glance

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}	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (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	363	<div> <div>25%</div> <div> <div></div> <div>71%</div> <div>22%</div> <div>8%</div> </div> </div>
1	B	363	<div> <div>17%</div> <div> <div></div> <div>70%</div> <div>18%</div> <div>12%</div> </div> </div>
1	C	363	<div> <div>13%</div> <div> <div></div> <div>79%</div> <div>15%</div> <div>6%</div> </div> </div>
1	D	363	<div> <div>12%</div> <div> <div></div> <div>78%</div> <div>14%</div> <div>8%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10753 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 Transforming growth factor beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	335	Total	C	N	O	S	0	0	0
			2684	1706	469	492	17			
1	B	320	Total	C	N	O	S	0	0	0
			2556	1620	452	467	17			
1	C	340	Total	C	N	O	S	0	0	0
			2728	1732	481	498	17			
1	D	335	Total	C	N	O	S	0	0	0
			2690	1711	473	489	17			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P07200
A	0	PRO	-	expression tag	UNP P07200
A	4	SER	CYS	engineered mutation	UNP P07200
A	85	VAL	LEU	variant	UNP P07200
A	147	GLN	ASN	engineered mutation	UNP P07200
B	-1	GLY	-	expression tag	UNP P07200
B	0	PRO	-	expression tag	UNP P07200
B	4	SER	CYS	engineered mutation	UNP P07200
B	85	VAL	LEU	variant	UNP P07200
B	147	GLN	ASN	engineered mutation	UNP P07200
C	-1	GLY	-	expression tag	UNP P07200
C	0	PRO	-	expression tag	UNP P07200
C	4	SER	CYS	engineered mutation	UNP P07200
C	85	VAL	LEU	variant	UNP P07200
C	147	GLN	ASN	engineered mutation	UNP P07200
D	-1	GLY	-	expression tag	UNP P07200
D	0	PRO	-	expression tag	UNP P07200
D	4	SER	CYS	engineered mutation	UNP P07200
D	85	VAL	LEU	variant	UNP P07200
D	147	GLN	ASN	engineered mutation	UNP P07200

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



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

- Molecule 3 is BETA-D-MANNOSE (three-letter code: BMA) (formula: $C_6H_{12}O_6$).

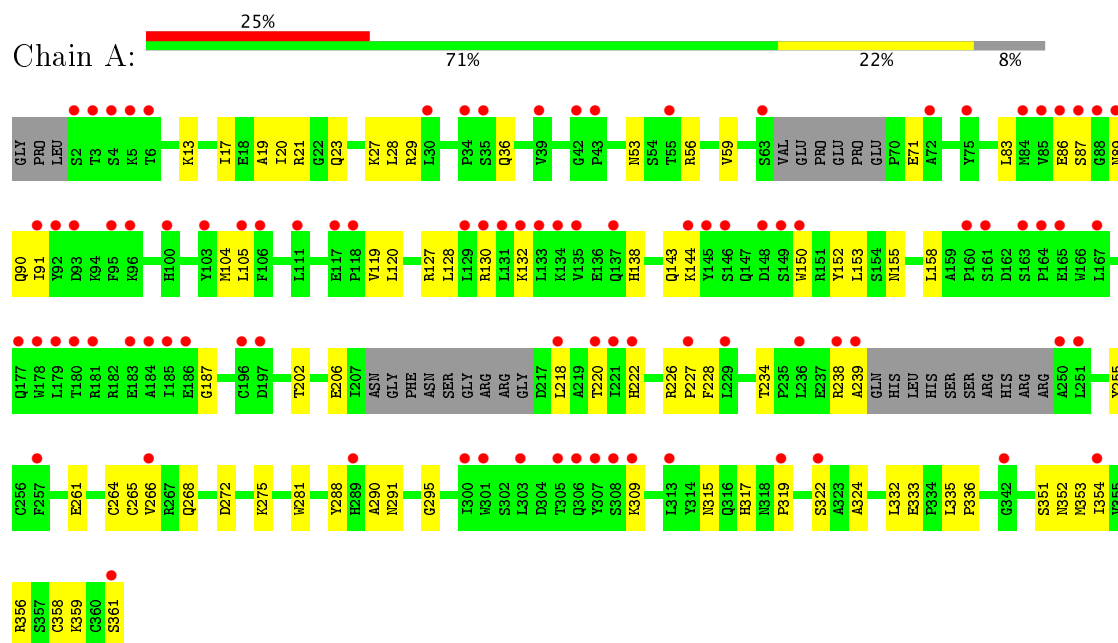


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			11	6	5		

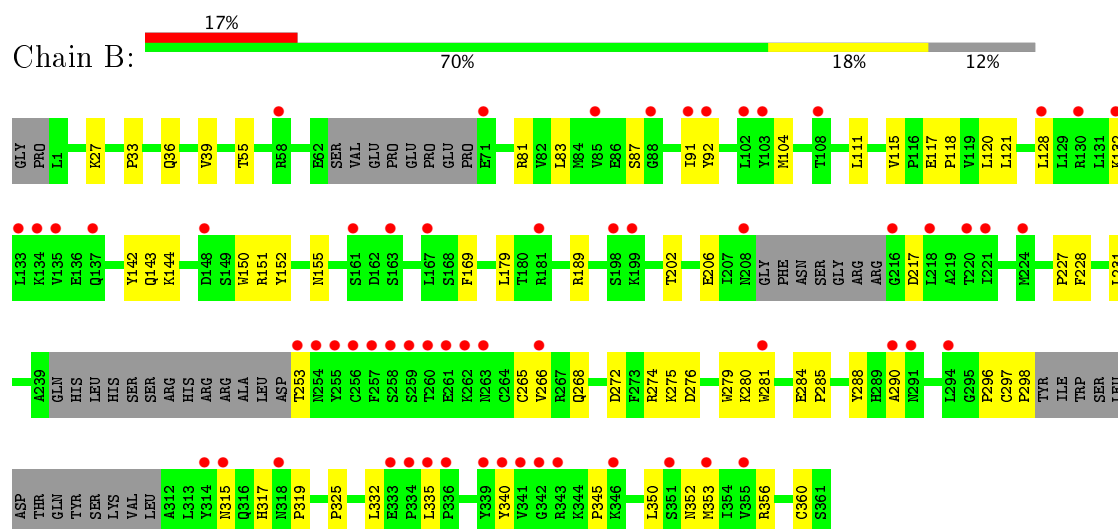
3 Residue-property plots

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

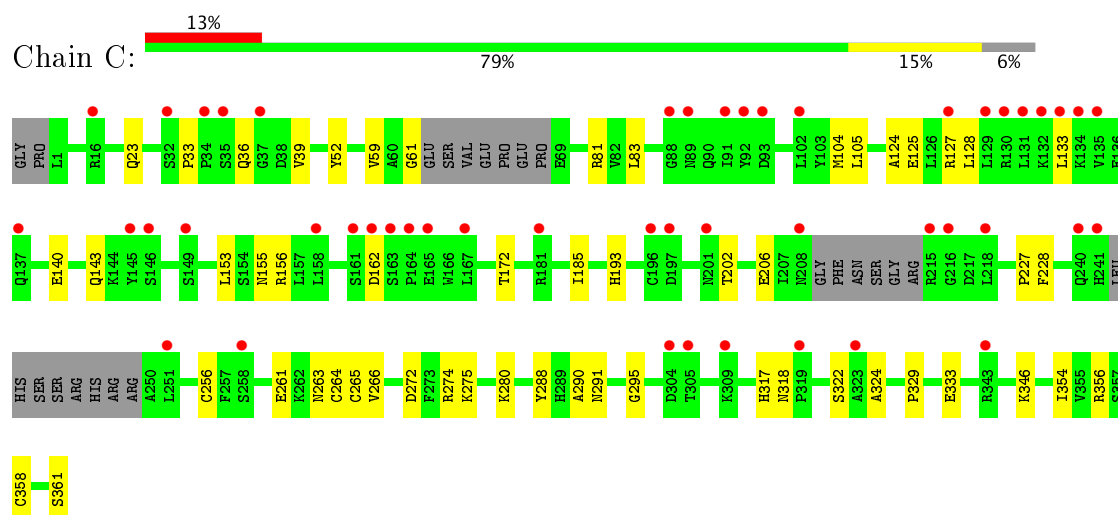
• Molecule 1: Transforming growth factor beta-1



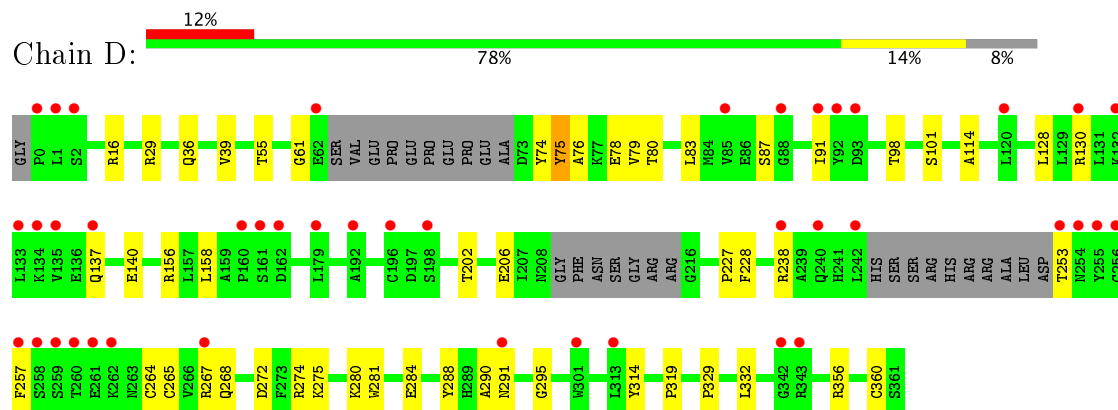
• Molecule 1: Transforming growth factor beta-1



• Molecule 1: Transforming growth factor beta-1



- Molecule 1: Transforming growth factor beta-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.66Å 126.92Å 137.92Å 90.00° 96.70° 90.00°	Depositor
Resolution (Å)	37.14 – 2.90 37.14 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.2 (37.14-2.90) 96.3 (37.14-2.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 2.90Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.239 , 0.280 0.238 , 0.279	Depositor DCC
R_{free} test set	1352 reflections (3.39%)	DCC
Wilson B-factor (Å ²)	129.9	Xtriage
Anisotropy	0.333	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 130.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10753	wwPDB-VP
Average B, all atoms (Å ²)	189.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.25% 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: BMA, 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	0.29	0/2747	0.49	1/3723 (0.0%)
1	B	0.28	0/2613	0.50	1/3537 (0.0%)
1	C	0.29	0/2792	0.50	1/3784 (0.0%)
1	D	0.33	0/2754	0.51	1/3732 (0.0%)
All	All	0.30	0/10906	0.50	4/14776 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	265	CYS	CA-CB-SG	9.06	130.31	114.00
1	A	265	CYS	CA-CB-SG	6.90	126.42	114.00
1	B	265	CYS	CA-CB-SG	6.30	125.34	114.00
1	C	265	CYS	CA-CB-SG	5.35	123.62	114.00

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	2684	0	2676	60	0
1	B	2556	0	2552	51	0
1	C	2728	0	2720	34	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2690	0	2687	30	0
2	B	28	0	24	2	0
2	C	28	0	25	3	0
2	D	28	0	25	1	0
3	B	11	0	10	0	0
All	All	10753	0	10719	158	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 7.

All (158) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:LEU:HD13	1:B:228:PHE:HB3	1.54	0.88
1:A:83:LEU:HD13	1:A:228:PHE:HB3	1.58	0.84
1:A:86:GLU:OE2	1:A:222:HIS:ND1	2.11	0.82
1:A:36:GLN:NE2	1:B:281:TRP:HB3	1.96	0.80
1:A:155:ASN:ND2	1:B:152:TYR:OH	2.16	0.79
1:D:83:LEU:HD13	1:D:228:PHE:HB3	1.64	0.78
1:B:298:PRO:HG2	1:B:317:HIS:HE1	1.51	0.75
1:D:288:TYR:CE2	1:D:290:ALA:HB2	2.23	0.74
1:D:29:ARG:HH11	1:D:75:TYR:HE1	1.36	0.72
1:C:288:TYR:CE2	1:C:290:ALA:HB2	2.24	0.72
1:B:132:LYS:NZ	1:B:217:ASP:O	2.24	0.71
1:B:143:GLN:OE1	1:B:151:ARG:NH2	2.25	0.69
1:B:298:PRO:HG2	1:B:317:HIS:CE1	2.28	0.68
1:A:255:TYR:CE2	1:A:261:GLU:OE2	2.49	0.65
1:C:83:LEU:HD13	1:C:228:PHE:HB3	1.78	0.64
1:A:288:TYR:CE2	1:A:290:ALA:HB2	2.32	0.64
1:A:272:ASP:HB3	1:A:275:LYS:HB3	1.79	0.64
1:A:23:GLN:HG2	1:A:27:LYS:NZ	2.13	0.63
1:C:128:LEU:HD23	1:C:227:PRO:HG3	1.82	0.62
1:A:36:GLN:HE21	1:B:281:TRP:HB3	1.65	0.61
1:D:264:CYS:HA	1:D:295:GLY:HA3	1.81	0.61
1:B:288:TYR:CE2	1:B:290:ALA:HB2	2.36	0.61
1:B:117:GLU:HB2	1:B:120:LEU:HD23	1.82	0.61
1:D:76:ALA:HB3	1:D:238:ARG:HH12	1.67	0.60
1:A:143:GLN:HB2	1:A:153:LEU:HD11	1.84	0.60
1:C:52:TYR:CE2	1:D:284:GLU:HB3	2.37	0.59
1:B:87:SER:HA	1:B:91:ILE:HD11	1.85	0.59
2:B:3053:NAG:H83	2:B:3053:NAG:O3	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:206:GLU:HG2	1:C:202:THR:HG22	1.84	0.58
1:C:105:LEU:HB3	1:C:185:ILE:HD11	1.84	0.58
1:D:253:THR:HG23	1:D:356:ARG:HD3	1.83	0.58
1:D:87:SER:HA	1:D:91:ILE:HD11	1.83	0.58
1:A:29:ARG:HD3	1:A:238:ARG:HB3	1.86	0.58
1:A:53:ASN:HD22	1:A:56:ARG:HH21	1.51	0.58
1:A:53:ASN:HD22	1:A:56:ARG:NH2	2.01	0.57
1:C:291:ASN:O	1:D:16:ARG:NH2	2.38	0.56
1:B:128:LEU:HD23	1:B:227:PRO:HG3	1.87	0.56
1:D:74:TYR:HD2	1:D:75:TYR:CD2	2.25	0.55
1:C:317:HIS:ND1	1:C:361:SER:O	2.38	0.55
1:A:23:GLN:HG2	1:A:27:LYS:HZ3	1.73	0.54
2:C:3053:NAG:O3	2:C:3053:NAG:H83	2.08	0.54
1:D:128:LEU:HD23	1:D:227:PRO:HG3	1.90	0.54
1:B:115:VAL:HG11	1:B:121:LEU:HB2	1.90	0.53
1:D:130:ARG:HH11	1:D:158:LEU:HB3	1.72	0.53
1:B:272:ASP:HB3	1:B:275:LYS:HB3	1.90	0.53
1:C:264:CYS:HA	1:C:295:GLY:HA3	1.90	0.53
2:C:3053:NAG:O3	2:C:3054:NAG:O5	2.24	0.53
1:C:333:GLU:HB2	1:C:354:ILE:HB	1.90	0.52
1:A:19:ALA:HB1	1:A:319:PRO:HB2	1.91	0.52
1:A:315:ASN:O	1:B:352:ASN:ND2	2.42	0.52
1:A:333:GLU:OE2	1:A:356:ARG:NE	2.41	0.52
1:C:59:VAL:HG11	1:C:127:ARG:NH1	2.25	0.51
1:C:125:GLU:OE1	1:C:127:ARG:NH2	2.43	0.51
1:A:13:LYS:NZ	1:B:268:GLN:O	2.44	0.51
1:B:315:ASN:O	1:B:319:PRO:HG3	2.11	0.51
1:A:119:VAL:HG13	1:A:120:LEU:HD22	1.92	0.50
1:A:120:LEU:HD11	1:A:239:ALA:HB2	1.92	0.50
1:A:130:ARG:HG3	1:A:218:LEU:HD11	1.94	0.50
1:B:128:LEU:HD12	1:B:169:PHE:HE2	1.76	0.50
1:A:264:CYS:HA	1:A:295:GLY:HA3	1.94	0.50
1:A:17:ILE:HD13	1:B:276:ASP:HB3	1.94	0.50
1:D:274:ARG:HH12	1:D:280:LYS:HD3	1.77	0.50
1:B:274:ARG:HH12	1:B:280:LYS:HE2	1.77	0.50
1:D:329:PRO:HB2	1:D:332:LEU:HD21	1.93	0.50
1:A:105:LEU:HG	1:A:187:GLY:HA3	1.94	0.49
1:A:120:LEU:HD12	1:A:234:THR:HB	1.93	0.49
1:A:351:SER:O	1:B:27:LYS:HE2	2.12	0.49
2:B:3053:NAG:O3	2:B:3054:NAG:O5	2.27	0.49
1:A:144:LYS:HB2	1:A:150:TRP:CZ3	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:GLU:HG2	1:D:202:THR:HG22	1.94	0.49
1:A:333:GLU:HB2	1:A:354:ILE:HB	1.95	0.49
1:B:297:CYS:O	1:B:325:PRO:HG2	2.13	0.48
1:B:202:THR:HG22	1:C:206:GLU:HG2	1.95	0.48
1:B:83:LEU:HA	1:B:228:PHE:HB3	1.95	0.48
1:C:143:GLN:HB2	1:C:153:LEU:HD11	1.94	0.48
1:A:255:TYR:OH	1:A:309:LYS:HD3	2.14	0.48
1:A:104:MET:SD	1:A:128:LEU:HD21	2.54	0.48
1:B:296:PRO:HB2	1:B:298:PRO:HD3	1.95	0.47
1:B:332:LEU:HD13	1:B:352:ASN:HB3	1.96	0.47
1:D:319:PRO:HD2	1:D:360:CYS:O	2.15	0.47
1:B:335:LEU:HB3	1:B:353:MET:HG3	1.96	0.47
1:B:144:LYS:HB2	1:B:150:TRP:CZ3	2.50	0.47
1:B:36:GLN:OE1	1:B:39:VAL:HG21	2.15	0.47
1:A:281:TRP:CD1	1:B:33:PRO:HB3	2.50	0.47
1:C:61:GLY:HA2	1:C:156:ARG:NH1	2.30	0.47
2:D:3053:NAG:H62	2:D:3054:NAG:C1	2.45	0.47
1:A:132:LYS:HZ3	1:A:218:LEU:HA	1.80	0.47
1:A:317:HIS:NE2	1:A:319:PRO:HG3	2.30	0.47
1:D:75:TYR:N	1:D:75:TYR:CD2	2.83	0.47
1:B:319:PRO:HD2	1:B:360:CYS:O	2.14	0.46
1:A:20:ILE:HA	1:A:23:GLN:HB3	1.98	0.46
1:A:202:THR:HG22	1:D:206:GLU:HG2	1.96	0.46
1:A:152:TYR:OH	1:B:155:ASN:ND2	2.47	0.46
1:A:27:LYS:HD2	1:B:350:LEU:HD23	1.98	0.46
1:C:261:GLU:HG3	1:C:263:ASN:H	1.80	0.46
1:A:83:LEU:HD12	1:A:226:ARG:HG2	1.97	0.45
1:D:257:PHE:CE1	1:D:268:GLN:HG3	2.51	0.45
1:A:87:SER:HA	1:A:91:ILE:HD11	1.99	0.45
1:C:256:CYS:SG	1:C:266:VAL:N	2.81	0.45
1:A:27:LYS:HE2	1:B:353:MET:SD	2.56	0.45
1:C:333:GLU:OE2	1:C:356:ARG:NE	2.39	0.45
1:C:124:ALA:HB3	1:C:172:THR:HA	1.98	0.45
1:B:274:ARG:NH1	1:B:280:LYS:HE2	2.32	0.45
1:A:336:PRO:HB3	1:B:55:THR:HG21	1.99	0.44
1:A:20:ILE:HG12	1:A:23:GLN:OE1	2.18	0.44
1:C:36:GLN:HA	1:C:39:VAL:HG23	2.00	0.44
1:A:255:TYR:CZ	1:A:261:GLU:OE2	2.70	0.44
1:A:220:THR:O	1:A:227:PRO:HD3	2.17	0.44
1:D:36:GLN:OE1	1:D:39:VAL:HG21	2.18	0.44
1:D:137:GLN:HG3	1:D:158:LEU:HD12	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:272:ASP:HB3	1:D:275:LYS:HB3	2.00	0.43
1:B:284:GLU:HA	1:B:285:PRO:HA	1.78	0.43
1:A:130:ARG:CZ	1:A:158:LEU:HB3	2.48	0.43
1:A:335:LEU:HB3	1:A:353:MET:HG3	2.00	0.43
1:A:128:LEU:HD23	1:A:227:PRO:HG3	2.00	0.43
1:A:322:SER:C	1:A:324:ALA:H	2.20	0.43
1:A:89:ASN:OD1	1:A:90:GLN:HG2	2.19	0.43
1:B:111:LEU:HD22	1:B:231:LEU:HD11	2.00	0.43
1:C:193:HIS:HE1	1:D:140:GLU:OE2	2.01	0.43
1:D:61:GLY:HA2	1:D:156:ARG:NH1	2.33	0.43
1:A:36:GLN:HE22	1:B:281:TRP:HB3	1.79	0.43
1:C:23:GLN:NE2	1:C:318:ASN:OD1	2.49	0.43
1:D:267:ARG:O	1:D:291:ASN:HB3	2.19	0.43
1:B:142:TYR:HB2	1:B:189:ARG:HB3	2.01	0.42
1:C:266:VAL:HG22	1:C:358:CYS:SG	2.59	0.42
1:C:346:LYS:NZ	1:D:78:GLU:OE2	2.51	0.42
1:A:130:ARG:NH1	1:A:158:LEU:HB3	2.34	0.42
1:B:118:PRO:HB3	1:B:179:LEU:HB2	2.01	0.42
1:B:340:TYR:CZ	1:B:345:PRO:HB3	2.55	0.42
1:B:91:ILE:HG13	1:B:92:TYR:CD1	2.55	0.42
1:C:272:ASP:HB3	1:C:275:LYS:HB3	2.02	0.42
1:B:104:MET:SD	1:B:128:LEU:HD21	2.59	0.42
1:B:81:ARG:HD2	1:B:228:PHE:CE2	2.55	0.41
1:D:98:THR:HB	1:D:101:SER:HB3	2.02	0.41
1:A:332:LEU:HD22	1:A:352:ASN:HB3	2.02	0.41
1:C:322:SER:C	1:C:324:ALA:H	2.24	0.41
1:C:81:ARG:HD2	1:C:228:PHE:CD2	2.55	0.41
1:D:55:THR:HG21	1:D:79:VAL:HG11	2.02	0.41
1:B:340:TYR:CE2	1:B:345:PRO:HB3	2.56	0.41
1:A:266:VAL:HG22	1:A:358:CYS:SG	2.61	0.41
1:C:127:ARG:HD2	1:C:228:PHE:CZ	2.55	0.41
1:C:274:ARG:HH12	1:C:280:LYS:HD3	1.85	0.41
1:C:33:PRO:HB3	1:D:281:TRP:CD1	2.55	0.41
2:C:3053:NAG:H83	2:C:3053:NAG:C3	2.51	0.41
1:C:329:PRO:HA	1:C:358:CYS:HA	2.02	0.41
1:A:359:LYS:HE2	1:A:361:SER:OG	2.21	0.41
1:B:253:THR:HG22	1:B:266:VAL:HG11	2.03	0.41
1:A:59:VAL:HG11	1:A:127:ARG:CZ	2.51	0.41
1:A:132:LYS:NZ	1:A:218:LEU:HA	2.36	0.41
1:B:132:LYS:HA	1:B:132:LYS:HD3	1.88	0.41
1:C:140:GLU:HG2	1:C:155:ASN:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:104:MET:SD	1:C:128:LEU:HD21	2.61	0.41
1:A:28:LEU:HD22	1:A:238:ARG:HH22	1.85	0.41
1:A:21:ARG:HG3	1:B:279:TRP:HZ2	1.85	0.40
1:C:133:LEU:HD13	1:C:162:ASP:O	2.22	0.40
1:A:268:GLN:HA	1:A:291:ASN:OD1	2.21	0.40
1:B:253:THR:HG23	1:B:356:ARG:HD3	2.04	0.40
1:D:80:THR:OG1	1:D:114:ALA:HB1	2.22	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	327/363 (90%)	318 (97%)	9 (3%)	0	100	100
1	B	310/363 (85%)	299 (96%)	11 (4%)	0	100	100
1	C	332/363 (92%)	320 (96%)	12 (4%)	0	100	100
1	D	327/363 (90%)	317 (97%)	10 (3%)	0	100	100
All	All	1296/1452 (89%)	1254 (97%)	42 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	299/323 (93%)	297 (99%)	2 (1%)	87	97
1	B	284/323 (88%)	284 (100%)	0	100	100
1	C	303/323 (94%)	303 (100%)	0	100	100
1	D	300/323 (93%)	298 (99%)	2 (1%)	87	97
All	All	1186/1292 (92%)	1182 (100%)	4 (0%)	94	98

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	GLU
1	A	138	HIS
1	D	75	TYR
1	D	314	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	155	ASN
1	B	317	HIS
1	D	263	ASN
1	D	316	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

7 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	B	3053	1,2	14,14,15	1.41	1 (7%)	15,19,21	3.01	4 (26%)
2	NAG	B	3054	3,2	14,14,15	0.69	1 (7%)	15,19,21	0.78	0
3	BMA	B	3055	2	11,11,12	0.89	0	13,15,17	0.97	1 (7%)
2	NAG	C	3053	1,2	14,14,15	0.49	0	15,19,21	2.87	2 (13%)
2	NAG	C	3054	2	14,14,15	0.54	0	15,19,21	0.66	0
2	NAG	D	3053	1,2	14,14,15	1.70	1 (7%)	15,19,21	1.05	1 (6%)
2	NAG	D	3054	2	14,14,15	1.14	1 (7%)	15,19,21	1.41	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
2	NAG	B	3053	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	3054	3,2	-	0/6/23/26	0/1/1/1
3	BMA	B	3055	2	-	0/2/19/22	0/1/1/1
2	NAG	C	3053	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	3054	2	-	0/6/23/26	0/1/1/1
2	NAG	D	3053	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	3054	2	-	0/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	3053	NAG	O5-C1	-6.20	1.33	1.43
2	B	3054	NAG	C1-C2	-2.21	1.49	1.52
2	D	3054	NAG	O5-C1	3.94	1.50	1.43
2	B	3053	NAG	O5-C1	4.39	1.50	1.43

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3054	NAG	C1-O5-C5	-2.83	108.27	112.17
3	B	3055	BMA	O2-C2-C3	-2.47	105.33	110.17
2	B	3053	NAG	O3-C3-C2	-2.28	104.51	109.39
2	D	3053	NAG	O4-C4-C5	2.45	115.47	109.28
2	D	3054	NAG	C3-C4-C5	3.30	116.03	110.22
2	B	3053	NAG	C1-C2-N2	3.88	117.12	110.49
2	B	3053	NAG	C1-O5-C5	5.72	120.05	112.17
2	C	3053	NAG	C1-C2-N2	5.83	120.45	110.49
2	B	3053	NAG	C2-N2-C7	8.06	134.71	122.94
2	C	3053	NAG	C2-N2-C7	8.37	135.15	122.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3053	NAG	2	0
2	B	3054	NAG	1	0
2	C	3053	NAG	3	0
2	C	3054	NAG	1	0
2	D	3053	NAG	1	0
2	D	3054	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	335/363 (92%)	1.30	92 (27%) 1 0	122, 208, 301, 367	0
1	B	320/363 (88%)	1.17	61 (19%) 1 1	130, 195, 272, 364	0
1	C	340/363 (93%)	0.92	48 (14%) 3 2	92, 161, 257, 344	0
1	D	335/363 (92%)	1.05	42 (12%) 4 3	106, 160, 258, 308	0
All	All	1330/1452 (91%)	1.11	243 (18%) 1 1	92, 183, 274, 367	0

All (243) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	242	LEU	16.2
1	A	91	ILE	10.3
1	B	342	GLY	9.9
1	B	260	THR	8.6
1	D	258	SER	8.5
1	D	253	THR	8.3
1	A	305	THR	8.2
1	D	93	ASP	7.3
1	B	92	TYR	7.0
1	D	255	TYR	7.0
1	C	146	SER	7.0
1	D	0	PRO	6.9
1	C	92	TYR	6.8
1	B	91	ILE	6.8
1	D	88	GLY	6.8
1	C	145	TYR	6.5
1	D	240	GLN	6.5
1	D	256	CYS	6.4
1	D	257	PHE	6.2
1	D	254	ASN	6.2
1	B	256	CYS	6.1

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Mol	Chain	Res	Type	RSRZ
1	B	255	TYR	6.0
1	D	161	SER	5.8
1	A	238	ARG	5.7
1	B	258	SER	5.7
1	D	301	TRP	5.7
1	A	179	LEU	5.6
1	A	307	TYR	5.5
1	B	220	THR	5.3
1	B	134	LYS	5.3
1	C	93	ASP	5.3
1	C	218	LEU	5.2
1	D	259	SER	5.2
1	A	129	LEU	5.1
1	A	2	SER	5.1
1	B	133	LEU	4.9
1	A	221	ILE	4.9
1	A	186	GLU	4.8
1	B	132	LYS	4.7
1	D	260	THR	4.7
1	A	306	GLN	4.4
1	C	216	GLY	4.4
1	A	134	LYS	4.4
1	A	105	LEU	4.3
1	A	342	GLY	4.3
1	B	314	TYR	4.3
1	A	117	GLU	4.2
1	A	181	ARG	4.2
1	B	259	SER	4.2
1	B	261	GLU	4.1
1	C	132	LYS	4.1
1	C	208	ASN	4.1
1	D	261	GLU	4.0
1	A	361	SER	4.0
1	A	250	ALA	4.0
1	C	149	SER	4.0
1	A	34	PRO	3.9
1	D	91	ILE	3.9
1	C	130	ARG	3.9
1	B	135	VAL	3.9
1	B	71	GLU	3.9
1	A	222	HIS	3.8
1	C	131	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	D	135	VAL	3.8
1	A	92	TYR	3.7
1	A	180	THR	3.7
1	A	239	ALA	3.7
1	D	132	LYS	3.7
1	B	340	TYR	3.7
1	B	181	ARG	3.7
1	B	318	ASN	3.7
1	A	165	GLU	3.6
1	A	184	ALA	3.6
1	A	251	LEU	3.6
1	C	240	GLN	3.6
1	B	130	ARG	3.6
1	C	88	GLY	3.6
1	B	161	SER	3.6
1	A	89	ASN	3.6
1	D	92	TYR	3.5
1	A	229	LEU	3.5
1	B	341	VAL	3.4
1	A	106	PHE	3.4
1	B	343	ARG	3.4
1	D	262	LYS	3.4
1	C	135	VAL	3.4
1	A	63	SER	3.3
1	B	334	PRO	3.3
1	D	134	LYS	3.3
1	B	257	PHE	3.3
1	C	162	ASP	3.3
1	C	181	ARG	3.3
1	C	37	GLY	3.2
1	A	103	TYR	3.2
1	A	39	VAL	3.2
1	D	238	ARG	3.2
1	A	88	GLY	3.2
1	A	145	TYR	3.2
1	A	319	PRO	3.2
1	D	1	LEU	3.2
1	B	346	LYS	3.1
1	B	85	VAL	3.1
1	B	355	VAL	3.1
1	A	86	GLU	3.1
1	C	161	SER	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	183	GLU	3.1
1	B	167	LEU	3.1
1	A	309	LYS	3.1
1	C	89	ASN	3.1
1	C	167	LEU	3.0
1	A	177	GLN	3.0
1	D	267	ARG	3.0
1	C	309	LYS	3.0
1	A	354	ILE	3.0
1	D	2	SER	3.0
1	B	199	LYS	3.0
1	C	134	LYS	3.0
1	B	262	LYS	2.9
1	C	163	SER	2.9
1	A	236	LEU	2.9
1	A	161	SER	2.9
1	B	148	ASP	2.9
1	A	87	SER	2.9
1	B	163	SER	2.9
1	C	305	THR	2.9
1	A	84	MET	2.9
1	A	96	LYS	2.9
1	C	323	ALA	2.9
1	B	58	ARG	2.9
1	A	132	LYS	2.9
1	D	342	GLY	2.9
1	B	291	ASN	2.9
1	D	160	PRO	2.8
1	B	198	SER	2.8
1	A	42	GLY	2.8
1	B	254	ASN	2.8
1	B	263	ASN	2.8
1	B	266	VAL	2.8
1	D	179	LEU	2.8
1	A	72	ALA	2.8
1	A	137	GLN	2.8
1	C	215	ARG	2.8
1	A	266	VAL	2.8
1	C	133	LEU	2.8
1	D	130	ARG	2.7
1	C	102	LEU	2.7
1	D	62	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	162	ASP	2.7
1	A	218	LEU	2.7
1	B	290	ALA	2.7
1	D	291	ASN	2.6
1	A	300	ILE	2.6
1	A	131	LEU	2.6
1	A	167	LEU	2.6
1	A	4	SER	2.6
1	C	137	GLN	2.6
1	B	336	PRO	2.6
1	C	127	ARG	2.6
1	B	294	LEU	2.6
1	A	85	VAL	2.6
1	A	178	TRP	2.6
1	A	303	LEU	2.6
1	A	322	SER	2.6
1	C	165	GLU	2.6
1	A	308	SER	2.6
1	A	6	THR	2.5
1	C	197	ASP	2.5
1	A	163	SER	2.5
1	C	241	HIS	2.5
1	C	258	SER	2.5
1	A	313	LEU	2.5
1	D	137	GLN	2.5
1	D	313	LEU	2.5
1	B	108	THR	2.5
1	A	3	THR	2.5
1	A	220	THR	2.5
1	D	85	VAL	2.4
1	D	343	ARG	2.4
1	A	75	TYR	2.4
1	A	135	VAL	2.4
1	C	196	CYS	2.4
1	C	343	ARG	2.4
1	B	218	LEU	2.4
1	C	201	ASN	2.4
1	A	196	CYS	2.4
1	B	339	TYR	2.4
1	B	137	GLN	2.4
1	A	197	ASP	2.4
1	B	315	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	133	LEU	2.4
1	B	221	ILE	2.4
1	A	95	PHE	2.3
1	A	148	ASP	2.3
1	C	158	LEU	2.3
1	B	216	GLY	2.3
1	D	192	ALA	2.3
1	B	335	LEU	2.3
1	A	43	PRO	2.3
1	C	304	ASP	2.3
1	A	130	ARG	2.3
1	A	146	SER	2.3
1	D	198	SER	2.3
1	A	227	PRO	2.2
1	B	103	TYR	2.2
1	A	35	SER	2.2
1	A	149	SER	2.2
1	A	289	HIS	2.2
1	A	5	LYS	2.2
1	C	91	ILE	2.2
1	C	35	SER	2.2
1	A	100	HIS	2.2
1	C	16	ARG	2.2
1	B	224	MET	2.2
1	C	129	LEU	2.2
1	A	164	PRO	2.2
1	D	133	LEU	2.1
1	A	118	PRO	2.1
1	A	160	PRO	2.1
1	D	196	CYS	2.1
1	A	30	LEU	2.1
1	C	319	PRO	2.1
1	A	55	THR	2.1
1	A	93	ASP	2.1
1	C	34	PRO	2.1
1	B	128	LEU	2.1
1	C	251	LEU	2.1
1	B	253	THR	2.1
1	B	88	GLY	2.1
1	B	351	SER	2.1
1	C	32	SER	2.1
1	D	120	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	301	TRP	2.0
1	B	281	TRP	2.0
1	B	353	MET	2.0
1	B	208	ASN	2.0
1	A	185	ILE	2.0
1	A	144	LYS	2.0
1	B	333	GLU	2.0
1	A	111	LEU	2.0
1	A	150	TRP	2.0
1	A	257	PHE	2.0
1	B	102	LEU	2.0
1	C	164	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAG	C	3054	14/15	0.73	0.33	-	248,281,331,396	0
2	NAG	B	3053	14/15	0.91	0.14	-	159,212,228,250	0
2	NAG	D	3053	14/15	0.84	0.14	-	173,207,246,264	0
3	BMA	B	3055	11/12	0.41	0.36	-	381,388,392,393	0
2	NAG	C	3053	14/15	0.91	0.15	-	159,210,239,242	0
2	NAG	B	3054	14/15	0.85	0.20	-	241,286,333,396	0
2	NAG	D	3054	14/15	0.71	0.32	-	245,281,331,396	0

6.5 Other polymers

There are no such residues in this entry.