



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:28 AM UTC

PDB ID : 5CMN / pdb_00005cmn
Title : FLRT3 LRR domain in complex with LPHN3 Olfactomedin domain
Authors : Lu, Y.; Salzman, G.; Arac, D.
Deposited on : 2015-07-17
Resolution : 3.60 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

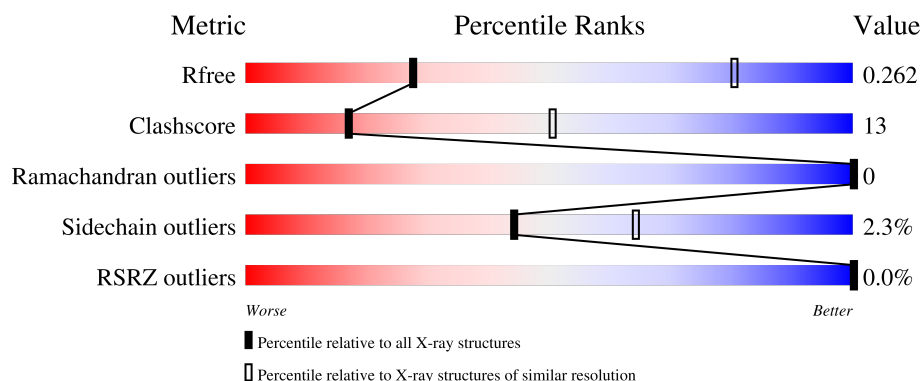
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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}	180053	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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 of 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	333	
1	B	333	
1	C	333	
1	D	333	
2	E	261	

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Mol	Chain	Length	Quality of chain
2	F	261	 68%25%• 5%
2	G	261	 69%25%• 5%
2	H	261	 70%23%• 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18420 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 Leucine-rich repeat transmembrane protein FLRT3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2575	1628	454	482	11			
1	B	321	Total	C	N	O	S	0	0	0
			2575	1628	454	482	11			
1	C	321	Total	C	N	O	S	0	0	0
			2575	1628	454	482	11			
1	D	321	Total	C	N	O	S	0	0	0
			2575	1628	454	482	11			

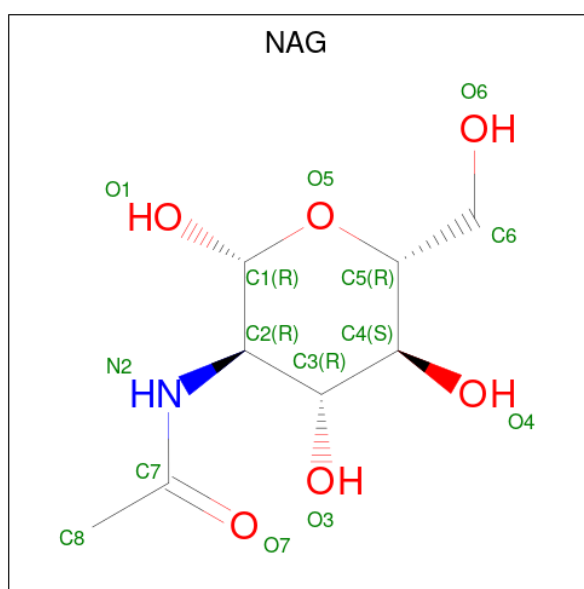
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	25	ALA	-	expression tag	UNP Q9NZU0
D	26	ASP	-	expression tag	UNP Q9NZU0
D	27	PRO	-	expression tag	UNP Q9NZU0
D	28	GLY	-	expression tag	UNP Q9NZU0
A	25	ALA	-	expression tag	UNP Q9NZU0
A	26	ASP	-	expression tag	UNP Q9NZU0
A	27	PRO	-	expression tag	UNP Q9NZU0
A	28	GLY	-	expression tag	UNP Q9NZU0
B	25	ALA	-	expression tag	UNP Q9NZU0
B	26	ASP	-	expression tag	UNP Q9NZU0
B	27	PRO	-	expression tag	UNP Q9NZU0
B	28	GLY	-	expression tag	UNP Q9NZU0
C	25	ALA	-	expression tag	UNP Q9NZU0
C	26	ASP	-	expression tag	UNP Q9NZU0
C	27	PRO	-	expression tag	UNP Q9NZU0
C	28	GLY	-	expression tag	UNP Q9NZU0

- Molecule 2 is a protein called Latrophilin-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	247	Total	C	N	O	S	0	0	0
			2015	1292	334	384	5			
2	F	247	Total	C	N	O	S	0	0	0
			2015	1292	334	384	5			
2	G	247	Total	C	N	O	S	0	0	0
			2015	1292	334	384	5			
2	H	247	Total	C	N	O	S	0	0	0
			2015	1292	334	384	5			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: 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		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	1	Total	Ca	0	0
			1	1		

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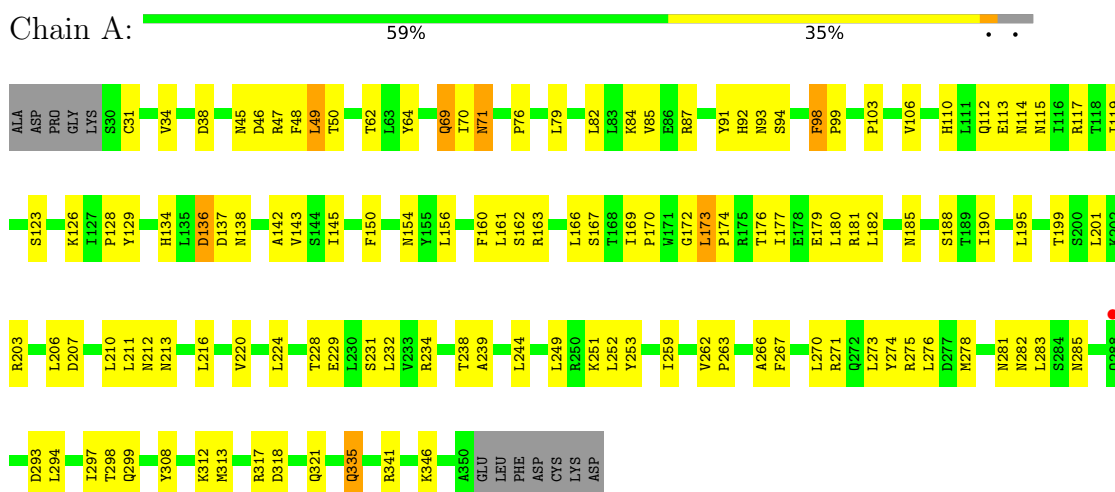
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	1	Total 1	Ca 1	0	0
4	G	1	Total 1	Ca 1	0	0
4	H	1	Total 1	Ca 1	0	0

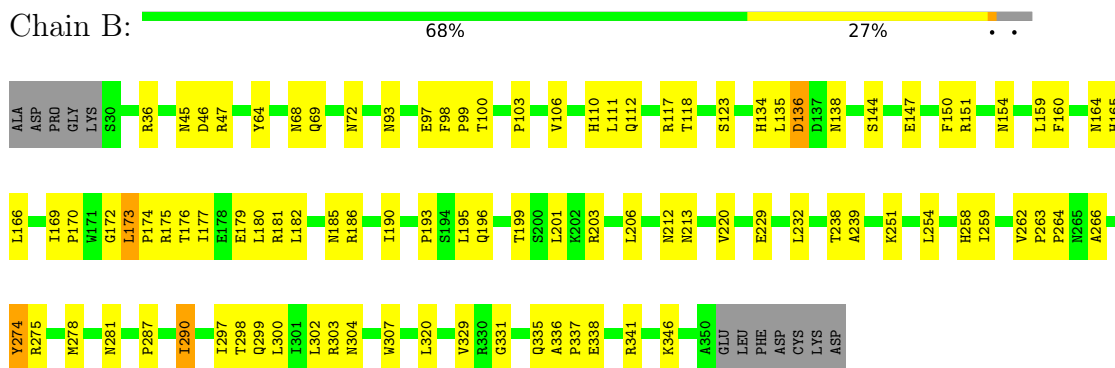
3 Residue-property plots

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.

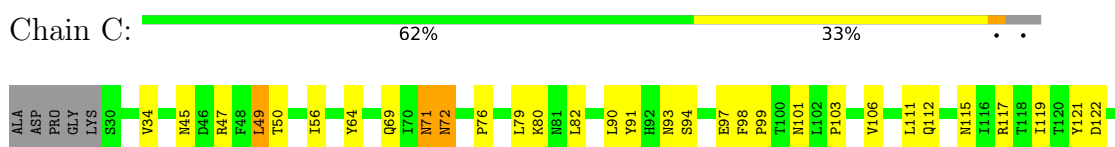
• Molecule 1: Leucine-rich repeat transmembrane protein FLRT3

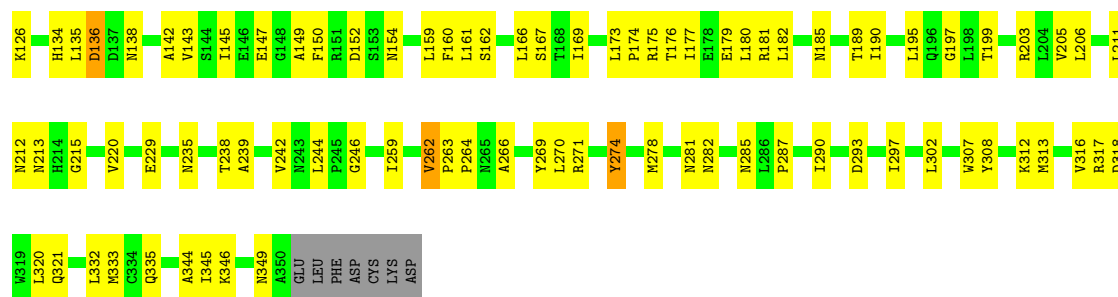


• Molecule 1: Leucine-rich repeat transmembrane protein FLRT3



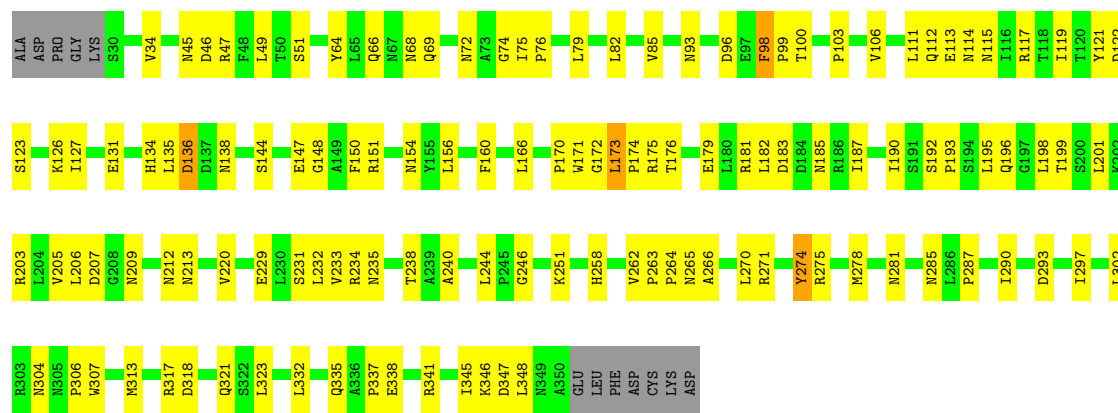
• Molecule 1: Leucine-rich repeat transmembrane protein FLRT3





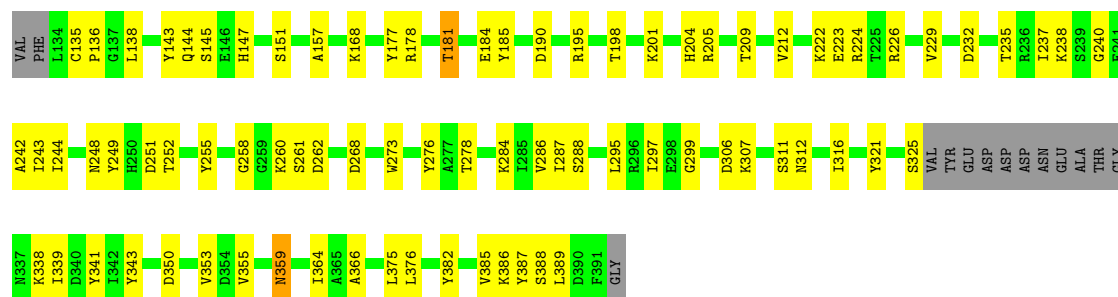
• Molecule 1: Leucine-rich repeat transmembrane protein FLRT3

Chain D: 59% 36%



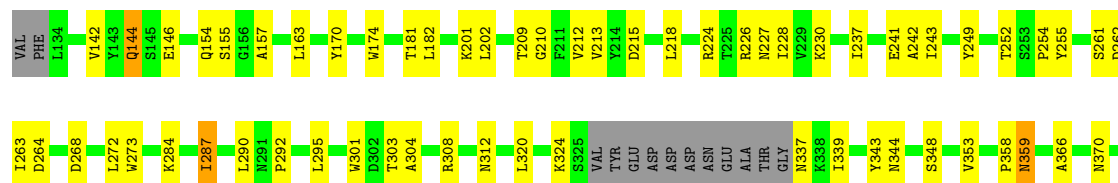
• Molecule 2: Latrophilin-3

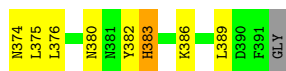
Chain E: 64% 30% 5%



• Molecule 2: Latrophilin-3

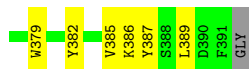
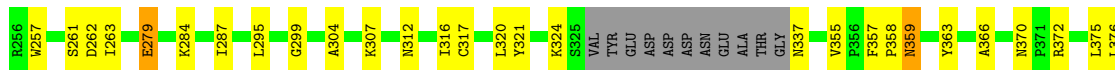
Chain F: 68% 25% 5%





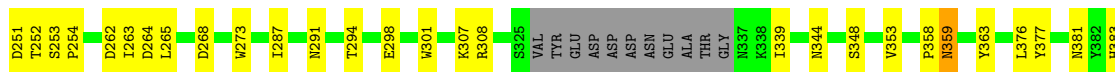
• Molecule 2: Latrophilin-3

Chain G: 69% 25% 5%



• Molecule 2: Latrophilin-3

Chain H: 70% 23% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	102.27Å 102.27Å 419.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	73.20 – 3.60 73.20 – 3.60	Depositor EDS
% Data completeness (in resolution range)	91.7 (73.20-3.60) 91.8 (73.20-3.60)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 3.58Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.195 , 0.262 0.197 , 0.262	Depositor DCC
R_{free} test set	1849 reflections (3.75%)	wwPDB-VP
Wilson B-factor (Å ²)	73.4	Xtriage
Anisotropy	0.382	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 93.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.350 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	18420	wwPDB-VP
Average B, all atoms (Å ²)	118.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.38% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.37	0/2627	0.86	5/3574 (0.1%)
1	B	0.36	0/2627	0.83	0/3574
1	C	0.38	0/2627	0.87	8/3574 (0.2%)
1	D	0.37	0/2627	0.83	3/3574 (0.1%)
2	E	0.39	0/2071	0.82	1/2820 (0.0%)
2	F	0.37	0/2071	0.81	2/2820 (0.1%)
2	G	0.38	0/2071	0.83	3/2820 (0.1%)
2	H	0.36	0/2071	0.79	1/2820 (0.0%)
All	All	0.37	0/18792	0.83	23/25576 (0.1%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	ILE	N-CA-C	6.43	114.90	107.89
1	C	169	ILE	CA-C-N	6.34	126.31	119.78
1	C	169	ILE	C-N-CA	6.34	126.31	119.78
2	G	355	VAL	CA-C-N	6.10	126.04	119.76
2	G	355	VAL	C-N-CA	6.10	126.04	119.76
1	C	50	THR	N-CA-C	-5.91	106.23	113.50
1	A	98	PHE	CA-C-N	5.90	126.36	119.99
1	A	98	PHE	C-N-CA	5.90	126.36	119.99
2	G	279	GLU	N-CA-C	-5.66	105.87	112.89
1	C	169	ILE	N-CA-C	5.57	113.96	107.89
2	H	229	VAL	N-CA-C	5.56	115.95	108.17
1	C	72	ASN	N-CA-C	5.51	119.26	112.54
1	D	98	PHE	CA-C-N	5.46	125.89	119.99
1	D	98	PHE	C-N-CA	5.46	125.89	119.99
2	F	202	LEU	CA-C-N	5.42	125.04	119.56
2	F	202	LEU	C-N-CA	5.42	125.04	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	75	ILE	N-CA-C	5.28	113.69	107.77
1	C	262	VAL	CA-C-N	5.23	125.76	120.38
1	C	262	VAL	C-N-CA	5.23	125.76	120.38
1	C	335	GLN	N-CA-C	-5.20	107.31	113.97
1	A	335	GLN	N-CA-C	-5.16	107.37	113.97
1	A	38	ASP	N-CA-C	5.04	115.64	107.23
2	E	311	SER	N-CA-C	-5.00	105.93	112.23

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	2575	0	2578	87	1
1	B	2575	0	2578	68	0
1	C	2575	0	2578	87	1
1	D	2575	0	2578	90	0
2	E	2015	0	1924	49	0
2	F	2015	0	1924	42	0
2	G	2015	0	1924	41	0
2	H	2015	0	1924	42	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
3	C	14	0	13	0	0
3	D	14	0	13	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
All	All	18420	0	18060	464	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (464) 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:D:264:PRO:HA	1:D:290:ILE:HG22	1.64	0.80
1:A:115:ASN:OD1	1:A:117:ARG:NH1	2.16	0.78
2:H:376:LEU:HB2	2:H:387:TYR:HB2	1.68	0.76
1:D:263:PRO:HG2	1:D:266:ALA:HB2	1.68	0.74
1:A:174:PRO:HG2	1:A:177:ILE:HG12	1.70	0.73
1:D:112:GLN:HA	1:D:138:ASN:HD21	1.52	0.73
1:C:115:ASN:OD1	1:C:117:ARG:NH1	2.23	0.72
1:B:251:LYS:HG2	1:B:275:ARG:HB3	1.71	0.70
1:A:251:LYS:HG2	1:A:275:ARG:HB3	1.73	0.70
2:H:287:ILE:HG13	2:H:301:TRP:HB2	1.73	0.70
1:C:271:ARG:NH2	1:C:293:ASP:OD2	2.26	0.68
2:E:242:ALA:HB2	2:E:295:LEU:HB3	1.74	0.68
2:E:316:ILE:HB	2:E:321:TYR:HE2	1.59	0.68
1:A:166:LEU:HB2	1:A:185:ASN:HD22	1.58	0.67
1:D:98:PHE:HB2	1:D:119:ILE:HG12	1.76	0.67
2:H:210:GLY:HA3	2:H:264:ASP:HA	1.77	0.67
2:H:198:THR:HG21	2:H:236:ARG:HH22	1.61	0.66
2:G:248:ASN:HB3	2:G:252:THR:HB	1.78	0.66
2:E:190:ASP:OD1	2:E:195:ARG:NH1	2.29	0.66
1:D:251:LYS:HG2	1:D:275:ARG:HB3	1.78	0.66
1:A:282:ASN:O	1:B:335:GLN:NE2	2.29	0.65
1:A:143:VAL:HG21	2:F:227:ASN:HD21	1.61	0.65
2:E:147:HIS:HA	2:E:385:VAL:HG12	1.77	0.65
1:C:263:PRO:HG2	1:C:266:ALA:HB2	1.78	0.65
2:F:375:LEU:HD13	2:F:386:LYS:HD3	1.77	0.65
1:A:134:HIS:HA	1:A:160:PHE:HB2	1.79	0.64
2:G:255:TYR:H	2:G:261:SER:HB2	1.63	0.63
1:A:136:ASP:O	1:A:138:ASN:ND2	2.32	0.63
2:F:210:GLY:HA3	2:F:264:ASP:HA	1.81	0.63
1:A:181:ARG:NH2	2:E:178:ARG:HG3	2.14	0.63
1:C:119:ILE:HD12	1:C:145:ILE:HG12	1.79	0.63
1:C:166:LEU:HB2	1:C:185:ASN:HD22	1.63	0.63
1:C:317:ARG:NH2	1:C:318:ASP:OD1	2.32	0.63
1:A:119:ILE:HD12	1:A:145:ILE:HG12	1.80	0.62
1:B:175:ARG:HD2	1:D:151:ARG:HH11	1.64	0.62
2:H:228:ILE:HD11	2:H:263:ILE:HG21	1.81	0.62
2:H:344:ASN:O	2:H:348:SER:N	2.32	0.62
1:A:45:ASN:ND2	1:A:46:ASP:OD1	2.32	0.62
1:A:166:LEU:HB2	1:A:185:ASN:ND2	2.13	0.62
2:E:284:LYS:HA	2:E:306:ASP:HA	1.81	0.62
1:D:264:PRO:HG2	2:E:299:GLY:HA3	1.80	0.62
1:B:110:HIS:HB3	1:B:112:GLN:HE22	1.65	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:117:ARG:NH2	1:D:113:GLU:HA	2.14	0.62
1:D:136:ASP:O	1:D:138:ASN:ND2	2.32	0.62
2:E:151:SER:N	2:E:382:TYR:O	2.31	0.62
1:B:100:THR:HG22	1:B:123:SER:HA	1.80	0.61
1:B:195:LEU:HD12	1:B:220:VAL:HG22	1.81	0.61
2:H:254:PRO:HA	2:H:262:ASP:HA	1.82	0.61
1:B:117:ARG:O	1:B:144:SER:HB3	1.99	0.61
2:G:376:LEU:HB2	2:G:387:TYR:HB2	1.82	0.61
1:C:285:ASN:HA	1:C:313:MET:HE1	1.83	0.61
2:E:204:HIS:HB3	2:E:222:LYS:HG3	1.82	0.61
1:C:76:PRO:HG2	1:C:79:LEU:HB2	1.81	0.61
2:G:161:ASP:HB3	2:G:164:GLN:HB2	1.82	0.61
2:G:376:LEU:HG	2:G:389:LEU:HD21	1.82	0.61
1:B:278:MET:O	1:B:281:ASN:ND2	2.34	0.60
1:C:34:VAL:HG12	1:C:47:ARG:HD3	1.83	0.60
1:D:181:ARG:NH2	1:D:203:ARG:HH12	1.99	0.60
1:B:179:GLU:HG3	1:B:203:ARG:HB3	1.83	0.60
1:C:189:THR:HG21	2:H:237:ILE:HA	1.83	0.60
1:D:117:ARG:O	1:D:144:SER:HB3	2.02	0.60
1:C:181:ARG:HH21	1:C:203:ARG:HH12	1.50	0.60
1:A:112:GLN:HA	1:A:138:ASN:HD21	1.66	0.60
1:C:112:GLN:HG3	2:G:224:ARG:HD3	1.84	0.60
1:C:166:LEU:HB2	1:C:185:ASN:ND2	2.17	0.59
1:A:271:ARG:NH2	1:A:293:ASP:OD2	2.35	0.59
1:B:182:LEU:HB2	1:B:206:LEU:HD23	1.84	0.59
1:D:181:ARG:HH21	1:D:203:ARG:HH12	1.48	0.59
1:B:212:ASN:OD1	1:B:213:ASN:N	2.36	0.59
2:H:229:VAL:HG22	2:H:241:GLU:HG3	1.84	0.59
1:C:72:ASN:ND2	1:C:97:GLU:HG2	2.18	0.59
1:B:151:ARG:HH11	1:D:175:ARG:HD2	1.68	0.58
2:H:248:ASN:HB3	2:H:252:THR:HB	1.85	0.58
1:D:166:LEU:HB2	1:D:185:ASN:ND2	2.18	0.58
1:A:206:LEU:HB2	1:A:232:LEU:HD23	1.86	0.58
1:C:302:LEU:HB3	1:C:307:TRP:HZ2	1.67	0.58
2:E:343:TYR:CD2	2:E:350:ASP:HB3	2.39	0.58
1:C:143:VAL:HG21	2:H:227:ASN:HD21	1.68	0.58
2:H:175:THR:O	2:H:179:THR:OG1	2.22	0.58
1:C:274:TYR:HA	1:C:297:ILE:HA	1.85	0.57
1:A:274:TYR:HA	1:A:297:ILE:HA	1.86	0.57
1:D:45:ASN:ND2	1:D:46:ASP:OD1	2.35	0.57
1:A:112:GLN:O	1:A:114:ASN:ND2	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:PRO:HG2	1:A:129:TYR:CD2	2.40	0.57
1:A:285:ASN:HA	1:A:313:MET:HE1	1.85	0.57
2:G:151:SER:N	2:G:382:TYR:O	2.32	0.57
1:B:166:LEU:HB2	1:B:185:ASN:HD22	1.70	0.57
1:D:302:LEU:HB3	1:D:307:TRP:HZ2	1.70	0.57
1:A:161:LEU:HB2	1:A:182:LEU:HD23	1.87	0.56
2:E:232:ASP:OD2	2:E:235:THR:OG1	2.17	0.56
1:B:274:TYR:HA	1:B:297:ILE:HA	1.88	0.56
1:D:321:GLN:NE2	1:D:346:LYS:O	2.39	0.56
1:D:212:ASN:OD1	1:D:213:ASN:N	2.34	0.56
1:D:278:MET:O	1:D:281:ASN:ND2	2.38	0.56
1:A:308:TYR:CZ	1:B:337:PRO:HB3	2.41	0.56
2:E:157:ALA:HB2	2:E:209:THR:HB	1.88	0.56
1:A:98:PHE:HD2	1:A:123:SER:HB3	1.70	0.55
1:C:190:ILE:N	1:C:215:GLY:O	2.34	0.55
1:D:166:LEU:HB2	1:D:185:ASN:HD22	1.69	0.55
2:F:268:ASP:OD1	2:F:273:TRP:NE1	2.40	0.55
2:H:206:VAL:HG13	2:H:220:PHE:HB2	1.89	0.55
2:G:316:ILE:HB	2:G:321:TYR:HE2	1.71	0.55
1:C:181:ARG:HH22	2:G:177:TYR:C	2.13	0.55
1:B:174:PRO:HG2	1:B:177:ILE:HG12	1.89	0.55
2:H:359:ASN:OD1	2:H:359:ASN:N	2.39	0.55
1:A:270:LEU:HD23	1:A:273:LEU:HD22	1.89	0.55
1:B:136:ASP:OD1	1:B:136:ASP:N	2.38	0.55
1:C:115:ASN:ND2	1:D:114:ASN:O	2.40	0.55
1:B:264:PRO:HG2	2:G:299:GLY:HA3	1.89	0.55
1:A:182:LEU:HB2	1:A:206:LEU:HD23	1.89	0.55
1:C:308:TYR:CZ	1:D:337:PRO:HB3	2.41	0.54
2:F:144:GLN:NE2	2:F:146:GLU:O	2.39	0.54
1:B:179:GLU:OE1	1:B:203:ARG:NH1	2.40	0.54
2:F:284:LYS:HD3	2:F:304:ALA:HA	1.89	0.54
1:C:278:MET:O	1:C:281:ASN:ND2	2.40	0.54
2:E:143:TYR:O	2:E:388:SER:N	2.40	0.54
1:A:150:PHE:HB3	1:A:174:PRO:HD3	1.89	0.54
1:D:111:LEU:HB2	1:D:135:LEU:HD23	1.89	0.54
1:D:346:LYS:HG3	1:D:347:ASP:OD2	2.08	0.54
2:E:268:ASP:OD1	2:E:273:TRP:NE1	2.39	0.54
1:B:181:ARG:HH21	1:B:203:ARG:HH12	1.56	0.54
2:G:359:ASN:N	2:G:359:ASN:OD1	2.41	0.54
1:B:229:GLU:HG2	1:B:251:LYS:HB2	1.90	0.54
1:D:244:LEU:HB2	1:D:270:LEU:HD11	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:51:SER:HA	1:D:74:GLY:HA3	1.89	0.54
1:B:36:ARG:HG2	1:B:47:ARG:HH21	1.73	0.53
1:D:190:ILE:HG22	1:D:220:VAL:HG21	1.89	0.53
2:E:255:TYR:H	2:E:261:SER:HB2	1.71	0.53
1:B:68:ASN:O	1:B:93:ASN:HA	2.08	0.53
1:B:103:PRO:HB2	1:B:106:VAL:HG23	1.90	0.53
1:C:332:LEU:HG	1:C:345:ILE:HD12	1.90	0.53
1:D:179:GLU:OE1	1:D:203:ARG:NH1	2.42	0.53
2:E:276:TYR:HA	2:E:307:LYS:HE3	1.90	0.53
2:E:376:LEU:HG	2:E:389:LEU:HD21	1.89	0.53
1:C:56:ILE:HG21	1:C:82:LEU:HD11	1.90	0.53
1:D:150:PHE:HB2	1:D:172:GLY:O	2.08	0.53
1:B:110:HIS:HB3	1:B:112:GLN:NE2	2.24	0.53
1:B:303:ARG:HA	1:B:331:GLY:HA3	1.91	0.52
1:B:263:PRO:HG2	1:B:266:ALA:HB2	1.91	0.52
1:C:94:SER:O	1:C:94:SER:OG	2.25	0.52
1:D:332:LEU:HG	1:D:345:ILE:HD12	1.90	0.52
1:D:76:PRO:HG2	1:D:79:LEU:HB2	1.91	0.52
1:D:238:THR:HA	1:D:258:HIS:O	2.09	0.52
2:E:376:LEU:HB2	2:E:387:TYR:HB2	1.91	0.52
2:G:242:ALA:HB2	2:G:295:LEU:HB3	1.90	0.52
1:C:47:ARG:NH1	2:G:279:GLU:OE2	2.43	0.52
2:E:205:ARG:HB2	2:E:223:GLU:HB2	1.91	0.52
1:A:231:SER:HA	1:A:253:TYR:HB2	1.91	0.52
1:C:134:HIS:HA	1:C:160:PHE:HB2	1.92	0.52
1:D:246:GLY:HA3	1:D:270:LEU:HA	1.91	0.52
1:C:122:ASP:O	1:C:126:LYS:HG2	2.10	0.52
1:C:181:ARG:HA	1:C:205:VAL:HB	1.92	0.52
1:D:156:LEU:HD23	1:D:174:PRO:HG3	1.92	0.52
2:E:229:VAL:HG13	2:E:238:LYS:HE2	1.92	0.52
1:B:147:GLU:HB2	1:D:196:GLN:NE2	2.25	0.52
1:B:170:PRO:HB2	1:B:173:LEU:HD13	1.92	0.52
1:C:82:LEU:O	1:C:103:PRO:HB3	2.10	0.52
2:G:228:ILE:HD11	2:G:263:ILE:HG21	1.91	0.52
1:D:317:ARG:HD2	1:D:348:LEU:O	2.09	0.52
1:C:112:GLN:HA	1:C:138:ASN:HD21	1.74	0.51
1:C:181:ARG:NH2	2:G:177:TYR:O	2.41	0.51
1:C:80:LYS:NZ	1:C:101:ASN:HA	2.26	0.51
1:C:239:ALA:HA	1:C:259:ILE:HG23	1.93	0.51
2:F:376:LEU:HG	2:F:389:LEU:HD21	1.93	0.51
1:A:113:GLU:H	1:A:137:ASP:HB2	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:GLN:HA	1:A:341:ARG:HG3	1.91	0.51
1:A:98:PHE:CD2	1:A:123:SER:HB3	2.46	0.50
1:B:264:PRO:HA	1:B:290:ILE:HG22	1.93	0.50
1:A:103:PRO:HB2	1:A:106:VAL:HG23	1.92	0.50
1:D:198:LEU:HB3	1:D:201:LEU:HB2	1.92	0.50
1:A:62:THR:HG21	1:A:87:ARG:HH21	1.77	0.50
2:E:359:ASN:N	2:E:359:ASN:OD1	2.42	0.50
1:A:163:ARG:HG2	1:B:186:ARG:NH2	2.27	0.50
2:H:157:ALA:HB2	2:H:209:THR:HB	1.94	0.50
1:C:181:ARG:NH2	1:C:203:ARG:HH12	2.10	0.50
1:C:203:ARG:HG3	1:C:229:GLU:HB2	1.93	0.50
1:C:321:GLN:NE2	1:C:346:LYS:O	2.45	0.50
1:D:317:ARG:NH2	1:D:318:ASP:OD1	2.44	0.50
2:G:147:HIS:HA	2:G:385:VAL:HG12	1.93	0.50
1:B:335:GLN:HA	1:B:341:ARG:HG3	1.93	0.50
1:D:72:ASN:HD21	1:D:96:ASP:H	1.60	0.50
2:E:278:THR:HG23	2:E:286:VAL:HG21	1.94	0.50
1:B:181:ARG:NH2	1:B:203:ARG:HH12	2.09	0.49
2:E:312:ASN:ND2	2:E:366:ALA:HA	2.26	0.49
1:C:179:GLU:HG3	1:C:203:ARG:HB3	1.93	0.49
1:C:212:ASN:OD1	1:C:213:ASN:N	2.45	0.49
1:D:34:VAL:HG12	1:D:47:ARG:HD3	1.93	0.49
1:D:103:PRO:O	1:D:106:VAL:HG23	2.13	0.49
2:E:258:GLY:O	2:E:261:SER:OG	2.25	0.49
1:A:212:ASN:OD1	1:A:213:ASN:N	2.44	0.49
1:C:161:LEU:HD13	1:C:166:LEU:HD11	1.95	0.49
2:E:226:ARG:HD3	2:E:244:ILE:O	2.12	0.49
1:C:211:LEU:HB2	1:C:235:ASN:OD1	2.12	0.49
2:F:272:LEU:HB3	2:F:290:LEU:O	2.12	0.49
1:D:100:THR:HG22	1:D:123:SER:HA	1.94	0.49
1:B:175:ARG:HD2	1:D:151:ARG:NH1	2.28	0.49
1:C:312:LYS:HG2	1:D:338:GLU:HG3	1.95	0.49
1:C:159:LEU:HB3	1:C:180:LEU:HD12	1.93	0.49
1:C:215:GLY:HA2	2:H:237:ILE:HG22	1.95	0.49
1:D:156:LEU:O	1:D:176:THR:OG1	2.28	0.49
2:E:375:LEU:HD13	2:E:386:LYS:HD3	1.95	0.49
2:G:172:MET:HE2	2:G:202:LEU:HD11	1.95	0.49
2:H:155:SER:HB2	2:H:174:TRP:HE3	1.77	0.49
1:A:91:TYR:HB3	1:A:112:GLN:HG2	1.95	0.49
1:A:244:LEU:HB2	1:A:270:LEU:HD11	1.95	0.48
2:F:255:TYR:H	2:F:261:SER:HB2	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:370:ASN:O	2:F:374:ASN:HA	2.13	0.48
2:H:291:ASN:HB2	2:H:298:GLU:CD	2.38	0.48
1:C:244:LEU:HB2	1:C:270:LEU:HD11	1.94	0.48
2:H:190:ASP:OD1	2:H:195:ARG:NH1	2.46	0.48
1:D:112:GLN:CD	1:D:112:GLN:H	2.21	0.48
1:C:182:LEU:HB2	1:C:206:LEU:HD23	1.96	0.48
1:A:161:LEU:HD13	1:A:166:LEU:HD11	1.96	0.48
1:C:64:TYR:CE1	2:G:252:THR:HA	2.49	0.48
2:H:291:ASN:OD1	2:H:294:THR:N	2.40	0.48
1:B:302:LEU:HB3	1:B:307:TRP:HZ2	1.79	0.47
1:D:240:ALA:HB3	1:D:263:PRO:HD2	1.95	0.47
2:F:324:LYS:O	2:F:337:ASN:HB2	2.14	0.47
1:A:203:ARG:HG3	1:A:229:GLU:HB2	1.97	0.47
2:F:287:ILE:HD13	2:F:320:LEU:HD22	1.96	0.47
1:D:182:LEU:HB2	1:D:206:LEU:HD23	1.96	0.47
1:A:49:LEU:O	1:A:69:GLN:N	2.42	0.47
1:A:207:ASP:HB3	1:A:234:ARG:NH1	2.29	0.47
2:E:181:THR:HG23	2:E:201:LYS:HD3	1.95	0.47
2:G:148:LEU:HB3	2:G:192:ILE:HG12	1.96	0.47
1:A:50:THR:HG22	1:A:69:GLN:HB2	1.96	0.47
1:D:68:ASN:O	1:D:93:ASN:HA	2.15	0.47
1:D:195:LEU:HB2	1:D:220:VAL:HG22	1.94	0.47
1:A:34:VAL:HG12	1:A:47:ARG:HD3	1.97	0.47
1:A:112:GLN:CD	1:A:112:GLN:H	2.23	0.47
1:A:239:ALA:HA	1:A:259:ILE:HG23	1.95	0.47
1:D:131:GLU:HA	1:D:156:LEU:HA	1.96	0.47
1:C:282:ASN:O	1:D:335:GLN:NE2	2.48	0.47
1:D:82:LEU:HB3	1:D:85:VAL:HG23	1.97	0.47
1:D:183:ASP:HB3	1:D:205:VAL:HG12	1.97	0.47
2:G:163:LEU:HD22	2:G:215:ASP:HB2	1.97	0.47
2:G:181:THR:HG22	2:G:201:LYS:HD3	1.97	0.47
1:A:298:THR:HG22	1:A:299:GLN:HG3	1.96	0.47
1:D:181:ARG:HA	1:D:205:VAL:HB	1.97	0.47
1:A:76:PRO:HG2	1:A:79:LEU:HB2	1.96	0.46
1:A:238:THR:HG22	1:B:304:ASN:ND2	2.30	0.46
1:D:45:ASN:HA	1:D:68:ASN:HD21	1.78	0.46
2:G:156:GLY:HA3	2:G:173:PRO:HA	1.97	0.46
2:G:287:ILE:HD13	2:G:320:LEU:HD22	1.97	0.46
1:D:274:TYR:HA	1:D:297:ILE:HA	1.97	0.46
1:C:238:THR:HG22	1:D:304:ASN:ND2	2.31	0.46
2:F:230:LYS:HD3	2:F:295:LEU:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:207:ASP:CG	2:H:223:GLU:HA	2.40	0.46
2:E:288:SER:HB3	2:E:297:ILE:HD12	1.97	0.46
2:F:312:ASN:ND2	2:F:366:ALA:HA	2.30	0.46
1:A:128:PRO:HG2	1:A:129:TYR:CE2	2.51	0.46
1:C:80:LYS:HZ3	1:C:101:ASN:HA	1.80	0.46
1:A:142:ALA:HB2	1:A:167:SER:OG	2.16	0.46
1:A:150:PHE:HB2	1:A:172:GLY:O	2.14	0.46
1:C:45:ASN:HB3	1:C:64:TYR:HB3	1.97	0.46
2:F:157:ALA:HB2	2:F:209:THR:HB	1.98	0.46
1:A:70:ILE:O	1:A:93:ASN:HB3	2.16	0.46
1:D:122:ASP:O	1:D:126:LYS:HG2	2.16	0.46
1:A:228:THR:O	1:A:249:LEU:HD12	2.16	0.46
1:B:136:ASP:O	1:B:138:ASN:ND2	2.49	0.46
1:B:182:LEU:HD12	1:B:206:LEU:HD21	1.97	0.46
1:C:121:TYR:HA	1:C:149:ALA:HA	1.98	0.46
2:E:276:TYR:HE2	2:E:278:THR:HG22	1.81	0.46
1:C:181:ARG:NH2	2:G:178:ARG:HG3	2.31	0.46
1:C:349:ASN:O	1:C:349:ASN:ND2	2.48	0.46
2:H:144:GLN:HG2	2:H:387:TYR:CE1	2.51	0.46
1:B:239:ALA:HA	1:B:259:ILE:HG23	1.98	0.45
1:C:136:ASP:HA	1:C:162:SER:O	2.15	0.45
2:F:228:ILE:HD11	2:F:263:ILE:HG21	1.98	0.45
1:A:136:ASP:OD1	2:E:224:ARG:NH2	2.49	0.45
1:B:232:LEU:HB2	1:B:254:LEU:HD23	1.97	0.45
2:E:248:ASN:HB3	2:E:252:THR:HB	1.98	0.45
2:F:252:THR:O	2:F:308:ARG:NH1	2.49	0.45
1:B:64:TYR:CE1	2:F:252:THR:HA	2.52	0.45
1:B:262:VAL:O	1:B:287:PRO:HG2	2.16	0.45
1:B:298:THR:HG22	1:B:299:GLN:HG3	1.98	0.45
2:H:251:ASP:N	2:H:262:ASP:OD1	2.46	0.45
2:F:182:LEU:HD22	2:F:218:LEU:HD22	1.97	0.45
2:E:226:ARG:HG3	2:E:243:ILE:HG23	1.99	0.45
2:E:238:LYS:HG2	2:E:240:GLY:H	1.82	0.45
1:A:136:ASP:HA	1:A:162:SER:O	2.16	0.45
1:B:196:GLN:NE2	1:D:147:GLU:HB2	2.32	0.45
2:F:230:LYS:NZ	2:F:292:PRO:O	2.35	0.45
2:F:249:TYR:HA	2:F:262:ASP:CB	2.46	0.45
1:D:115:ASN:OD1	1:D:117:ARG:NH1	2.50	0.45
2:F:226:ARG:HG3	2:F:243:ILE:HG23	1.98	0.45
1:B:169:ILE:HD12	1:B:190:ILE:HG23	1.98	0.45
1:C:175:ARG:HG2	1:C:197:GLY:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:121:TYR:HD1	1:D:148:GLY:O	2.00	0.45
2:F:181:THR:HG22	2:F:201:LYS:HD3	1.98	0.45
1:D:229:GLU:HG2	1:D:251:LYS:HB2	1.99	0.45
1:A:312:LYS:CB	1:B:338:GLU:HG3	2.47	0.44
1:B:112:GLN:H	1:B:112:GLN:CD	2.25	0.44
1:D:192:SER:HB2	1:D:193:PRO:HD3	1.99	0.44
1:D:207:ASP:HB3	1:D:234:ARG:NH1	2.33	0.44
1:D:209:ASN:O	1:D:235:ASN:HA	2.18	0.44
2:E:339:ILE:O	2:E:353:VAL:HG23	2.17	0.44
2:F:155:SER:HB2	2:F:174:TRP:CE3	2.52	0.44
2:F:255:TYR:CE2	2:F:312:ASN:HA	2.52	0.44
2:H:182:LEU:HD22	2:H:218:LEU:HD22	1.99	0.44
1:D:166:LEU:HB3	1:D:187:ILE:HD11	1.99	0.44
2:E:338:LYS:HD3	2:E:355:VAL:O	2.17	0.44
1:B:154:ASN:O	1:B:176:THR:HG21	2.17	0.44
1:D:134:HIS:HA	1:D:160:PHE:HB2	2.00	0.44
2:H:376:LEU:HG	2:H:389:LEU:HD21	1.99	0.44
1:A:154:ASN:O	1:A:176:THR:HG21	2.17	0.44
1:B:164:ASN:HB3	1:B:165:HIS:H	1.70	0.44
1:B:111:LEU:HB2	1:B:135:LEU:HD23	1.99	0.44
1:C:154:ASN:O	1:C:176:THR:HG21	2.17	0.44
2:F:163:LEU:HD22	2:F:215:ASP:HB2	1.99	0.44
2:G:136:PRO:HD2	2:G:317:CYS:SG	2.58	0.44
2:H:254:PRO:HD2	2:H:308:ARG:NH1	2.32	0.44
1:A:71:ASN:HB3	1:A:94:SER:OG	2.17	0.44
1:A:321:GLN:NE2	1:A:346:LYS:O	2.51	0.44
1:A:98:PHE:HA	1:A:99:PRO:HD3	1.69	0.44
1:A:317:ARG:NH2	1:A:318:ASP:OD1	2.51	0.44
1:B:45:ASN:O	1:B:47:ARG:NH1	2.51	0.44
1:B:193:PRO:HB2	1:D:171:TRP:CE3	2.53	0.44
1:C:143:VAL:HG21	2:H:227:ASN:ND2	2.31	0.44
1:D:287:PRO:HD2	1:D:290:ILE:HD13	2.00	0.44
1:D:335:GLN:HA	1:D:341:ARG:HG3	2.00	0.44
2:F:242:ALA:HB2	2:F:295:LEU:HB3	2.00	0.44
1:B:159:LEU:HB3	1:B:180:LEU:HD12	2.00	0.44
1:C:111:LEU:HB2	1:C:135:LEU:HD23	1.99	0.44
2:E:177:TYR:HA	2:E:223:GLU:OE2	2.17	0.44
2:H:155:SER:HB2	2:H:174:TRP:CE3	2.53	0.44
2:H:268:ASP:OD1	2:H:273:TRP:NE1	2.51	0.44
1:C:161:LEU:HB2	1:C:182:LEU:HD23	2.00	0.43
2:H:377:TYR:CE2	2:H:386:LYS:HE2	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:ILE:HG22	1:B:201:LEU:HD13	2.00	0.43
2:F:359:ASN:OD1	2:F:359:ASN:N	2.50	0.43
2:G:153:HIS:O	2:G:382:TYR:HA	2.18	0.43
2:G:284:LYS:HD3	2:G:304:ALA:HA	1.99	0.43
2:E:316:ILE:HB	2:E:321:TYR:CE2	2.45	0.43
1:A:195:LEU:HD12	1:A:220:VAL:HG22	2.00	0.43
1:C:264:PRO:HA	1:C:290:ILE:HG22	2.00	0.43
1:A:267:PHE:HB3	1:A:294:LEU:HD21	1.99	0.43
1:D:64:TYR:CE1	2:H:252:THR:HA	2.53	0.43
2:G:257:TRP:HB2	2:G:363:TYR:CZ	2.53	0.43
2:E:184:GLU:HB3	2:E:198:THR:HG22	2.01	0.43
2:E:325:SER:OG	2:E:364:ILE:HD12	2.19	0.43
1:A:190:ILE:HD11	1:A:211:LEU:HD11	1.99	0.43
2:G:170:TYR:HB3	2:G:182:LEU:HD11	2.01	0.43
1:A:64:TYR:CE1	2:E:252:THR:HA	2.53	0.43
1:A:259:ILE:HG22	1:A:283:LEU:HD21	2.00	0.43
1:C:103:PRO:O	1:C:106:VAL:HG23	2.18	0.43
1:C:271:ARG:HG2	1:C:293:ASP:O	2.19	0.43
2:H:228:ILE:HD12	2:H:265:LEU:HD11	2.01	0.43
1:B:45:ASN:ND2	1:B:46:ASP:OD1	2.44	0.43
1:B:238:THR:HA	1:B:258:HIS:O	2.19	0.43
1:C:142:ALA:HB2	1:C:167:SER:OG	2.19	0.43
1:C:195:LEU:HD12	1:C:220:VAL:HG22	2.00	0.43
2:F:226:ARG:HB3	2:F:249:TYR:CD1	2.54	0.43
2:H:253:SER:O	2:H:307:LYS:HD3	2.18	0.43
1:A:84:LYS:HA	1:A:84:LYS:HD3	1.80	0.42
1:A:94:SER:OG	1:A:94:SER:O	2.27	0.42
1:A:263:PRO:HG2	1:A:266:ALA:HB2	2.01	0.42
1:D:265:ASN:ND2	2:E:297:ILE:O	2.52	0.42
1:A:179:GLU:HG2	1:A:180:LEU:N	2.34	0.42
2:E:276:TYR:CE2	2:E:278:THR:HG22	2.54	0.42
2:G:312:ASN:ND2	2:G:366:ALA:HA	2.33	0.42
1:C:71:ASN:HB3	1:C:94:SER:O	2.20	0.42
2:H:339:ILE:O	2:H:353:VAL:HG23	2.19	0.42
1:A:252:LEU:HD23	1:A:276:LEU:HD13	2.02	0.42
1:B:134:HIS:HA	1:B:160:PHE:HB2	2.00	0.42
1:A:92:HIS:HA	1:A:113:GLU:O	2.20	0.42
1:B:72:ASN:ND2	1:B:97:GLU:HG2	2.34	0.42
1:D:154:ASN:O	1:D:176:THR:HG21	2.19	0.42
1:D:262:VAL:O	1:D:287:PRO:HG2	2.19	0.42
2:F:170:TYR:CE2	2:F:213:VAL:HG11	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:PHE:HA	1:A:69:GLN:OE1	2.19	0.42
1:B:110:HIS:CE1	1:B:134:HIS:CD2	3.07	0.42
1:C:112:GLN:HA	1:C:136:ASP:O	2.20	0.42
1:C:150:PHE:HB3	1:C:174:PRO:HD3	2.01	0.42
1:D:150:PHE:HB3	1:D:174:PRO:HD3	2.02	0.42
1:D:271:ARG:NH2	1:D:293:ASP:OD2	2.53	0.42
2:F:339:ILE:O	2:F:353:VAL:HG23	2.20	0.42
1:A:278:MET:O	1:A:281:ASN:ND2	2.52	0.42
1:B:112:GLN:HG3	2:F:224:ARG:HD3	2.02	0.42
1:B:151:ARG:NH1	1:D:175:ARG:HD2	2.33	0.42
1:C:317:ARG:O	1:C:321:GLN:HG2	2.19	0.42
1:D:170:PRO:HB2	1:D:173:LEU:HD13	2.01	0.42
2:E:144:GLN:HG3	2:E:145:SER:N	2.33	0.42
1:A:112:GLN:HA	1:A:136:ASP:O	2.20	0.42
1:A:308:TYR:HB2	1:B:336:ALA:HB1	2.00	0.42
1:B:98:PHE:HA	1:B:99:PRO:HD3	1.80	0.42
2:G:226:ARG:HD3	2:G:244:ILE:O	2.19	0.42
2:G:261:SER:O	2:G:307:LYS:NZ	2.38	0.42
2:G:324:LYS:O	2:G:337:ASN:HB2	2.20	0.42
1:A:91:TYR:HE1	2:E:249:TYR:CE2	2.37	0.41
1:A:110:HIS:HB3	1:A:112:GLN:NE2	2.35	0.41
1:A:170:PRO:HB2	1:A:173:LEU:HD13	2.01	0.41
1:B:300:LEU:HB3	1:B:329:VAL:HG22	2.02	0.41
2:F:344:ASN:O	2:F:348:SER:N	2.53	0.41
2:G:316:ILE:HB	2:G:321:TYR:CE2	2.53	0.41
1:A:201:LEU:HD23	1:A:224:LEU:HD13	2.02	0.41
1:C:181:ARG:HG2	1:C:205:VAL:HG21	2.02	0.41
1:C:246:GLY:HA3	1:C:269:TYR:O	2.19	0.41
1:D:98:PHE:HA	1:D:99:PRO:HD3	1.69	0.41
2:F:303:THR:HG22	2:F:343:TYR:CD2	2.56	0.41
1:C:242:VAL:HG12	1:C:266:ALA:HA	2.02	0.41
1:C:316:VAL:O	1:C:320:LEU:HG	2.21	0.41
2:F:230:LYS:HB2	2:F:295:LEU:HD21	2.02	0.41
2:F:254:PRO:HD2	2:F:308:ARG:NH1	2.36	0.41
2:H:207:ASP:N	2:H:221:ASN:O	2.43	0.41
1:A:182:LEU:HD12	1:A:206:LEU:HD21	2.02	0.41
1:A:188:SER:HA	1:A:210:LEU:O	2.20	0.41
1:D:126:LYS:HA	1:D:126:LYS:HD3	1.80	0.41
2:G:370:ASN:OD1	2:G:372:ARG:HB3	2.21	0.41
2:H:226:ARG:HD3	2:H:244:ILE:O	2.19	0.41
1:A:212:ASN:O	1:A:216:LEU:HD13	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:PHE:HA	1:C:99:PRO:HD3	1.83	0.41
1:C:282:ASN:ND2	1:D:306:PRO:HD3	2.35	0.41
1:D:205:VAL:HA	1:D:231:SER:HB3	2.02	0.41
2:E:251:ASP:H	2:E:262:ASP:CG	2.29	0.41
1:A:150:PHE:CB	1:A:174:PRO:HD3	2.50	0.41
1:C:136:ASP:O	1:C:138:ASN:ND2	2.52	0.41
1:D:45:ASN:HB2	1:D:66:GLN:HG2	2.03	0.41
1:D:263:PRO:HA	1:D:264:PRO:HD2	1.81	0.41
2:E:135:CYS:HA	2:E:136:PRO:HD3	1.89	0.41
1:C:174:PRO:HG2	1:C:177:ILE:HG12	2.02	0.41
1:A:262:VAL:HA	1:A:263:PRO:HD2	1.93	0.41
1:C:150:PHE:C	1:C:152:ASP:H	2.28	0.41
1:C:262:VAL:O	1:C:287:PRO:HG2	2.21	0.41
1:A:251:LYS:HD2	1:A:253:TYR:OH	2.21	0.41
1:B:112:GLN:HA	1:B:138:ASN:HD21	1.85	0.41
1:B:320:LEU:HD13	1:B:346:LYS:HB3	2.02	0.41
1:C:166:LEU:HD12	1:C:185:ASN:ND2	2.36	0.41
1:C:190:ILE:HD11	1:C:211:LEU:HD11	2.03	0.41
2:F:154:GLN:HB2	2:F:382:TYR:CZ	2.56	0.41
2:F:210:GLY:CA	2:F:264:ASP:HA	2.48	0.41
2:G:134:LEU:HD23	2:G:134:LEU:HA	1.93	0.41
2:G:170:TYR:CD2	2:G:213:VAL:HG21	2.56	0.41
2:G:357:PHE:HA	2:G:358:PRO:HD3	1.90	0.41
2:H:150:GLU:HG3	2:H:383:HIS:CD2	2.55	0.41
2:H:156:GLY:HA3	2:H:173:PRO:HA	2.02	0.41
1:C:112:GLN:H	1:C:112:GLN:CD	2.29	0.41
1:D:135:LEU:O	1:D:138:ASN:ND2	2.54	0.41
1:D:285:ASN:HA	1:D:313:MET:HE1	2.03	0.41
2:F:142:VAL:O	2:F:358:PRO:HD3	2.21	0.41
2:F:241:GLU:N	2:F:295:LEU:HD13	2.36	0.41
2:F:287:ILE:HG13	2:F:301:TRP:HB2	2.03	0.41
1:A:82:LEU:HB3	1:A:85:VAL:HG23	2.02	0.40
1:B:118:THR:HA	1:B:144:SER:O	2.20	0.40
1:C:91:TYR:HB2	1:C:112:GLN:HG2	2.03	0.40
1:D:64:TYR:OH	2:H:251:ASP:O	2.30	0.40
2:E:260:LYS:C	2:E:262:ASP:H	2.28	0.40
2:H:142:VAL:O	2:H:358:PRO:HD3	2.21	0.40
2:H:203:PRO:HG3	2:H:238:LYS:HB2	2.01	0.40
1:B:150:PHE:HB2	1:B:172:GLY:O	2.22	0.40
1:D:207:ASP:HA	1:D:233:VAL:O	2.21	0.40
2:G:375:LEU:HD13	2:G:386:LYS:HD3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:90:LEU:O	1:C:93:ASN:ND2	2.52	0.40
1:C:333:MET:HG2	1:C:344:ALA:HA	2.03	0.40
2:E:168:LYS:HA	2:E:185:TYR:O	2.22	0.40
2:E:341:TYR:OH	2:E:350:ASP:OD2	2.36	0.40
2:G:154:GLN:O	2:G:379:TRP:NE1	2.53	0.40
1:D:206:LEU:HB2	1:D:232:LEU:HD23	2.03	0.40
2:F:380:ASN:O	2:F:383:HIS:HB2	2.20	0.40
1:A:126:LYS:HA	1:A:126:LYS:HD3	1.88	0.40
1:A:156:LEU:HD12	1:A:156:LEU:HA	1.93	0.40
1:C:47:ARG:HB2	1:C:49:LEU:HD13	2.04	0.40
2:G:251:ASP:N	2:G:262:ASP:OD1	2.52	0.40
2:H:363:TYR:HB2	2:H:381:ASN:ND2	2.37	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:TYR:OH	1:C:147:GLU:O[4_565]	2.12	0.08

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	319/333 (96%)	296 (93%)	23 (7%)	0	100	100
1	B	319/333 (96%)	298 (93%)	21 (7%)	0	100	100
1	C	319/333 (96%)	295 (92%)	24 (8%)	0	100	100
1	D	319/333 (96%)	294 (92%)	25 (8%)	0	100	100
2	E	243/261 (93%)	229 (94%)	14 (6%)	0	100	100
2	F	243/261 (93%)	228 (94%)	15 (6%)	0	100	100
2	G	243/261 (93%)	227 (93%)	16 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	243/261 (93%)	228 (94%)	15 (6%)	0	100	100
All	All	2248/2376 (95%)	2095 (93%)	153 (7%)	0	100	100

There are no Ramachandran outliers to report.

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	296/306 (97%)	289 (98%)	7 (2%)	43	63
1	B	296/306 (97%)	290 (98%)	6 (2%)	48	66
1	C	296/306 (97%)	289 (98%)	7 (2%)	43	63
1	D	296/306 (97%)	288 (97%)	8 (3%)	39	61
2	E	219/230 (95%)	213 (97%)	6 (3%)	39	61
2	F	219/230 (95%)	213 (97%)	6 (3%)	39	61
2	G	219/230 (95%)	216 (99%)	3 (1%)	59	71
2	H	219/230 (95%)	215 (98%)	4 (2%)	51	68
All	All	2060/2144 (96%)	2013 (98%)	47 (2%)	44	64

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	CYS
1	A	49	LEU
1	A	69	GLN
1	A	71	ASN
1	A	136	ASP
1	A	173	LEU
1	A	199	THR
1	B	69	GLN
1	B	136	ASP
1	B	173	LEU

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Mol	Chain	Res	Type
1	B	199	THR
1	B	274	TYR
1	B	290	ILE
1	C	49	LEU
1	C	69	GLN
1	C	71	ASN
1	C	136	ASP
1	C	173	LEU
1	C	199	THR
1	C	274	TYR
1	D	49	LEU
1	D	69	GLN
1	D	127	ILE
1	D	136	ASP
1	D	173	LEU
1	D	199	THR
1	D	274	TYR
1	D	323	LEU
2	E	138	LEU
2	E	181	THR
2	E	212	VAL
2	E	237	ILE
2	E	287	ILE
2	E	359	ASN
2	F	144	GLN
2	F	212	VAL
2	F	237	ILE
2	F	287	ILE
2	F	359	ASN
2	F	383	HIS
2	G	138	LEU
2	G	212	VAL
2	G	359	ASN
2	H	144	GLN
2	H	179	THR
2	H	212	VAL
2	H	359	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	69	GLN

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Mol	Chain	Res	Type
1	B	134	HIS
1	B	282	ASN
1	B	296	ASN
1	C	213	ASN
1	C	282	ASN
1	C	285	ASN
1	C	296	ASN
1	D	72	ASN
1	D	138	ASN
1	D	213	ASN
1	D	285	ASN
1	D	349	ASN
2	E	344	ASN
2	G	154	GLN
2	G	204	HIS
2	H	154	GLN
2	H	227	ASN
2	H	246	ASN
2	H	270	ASN
2	H	381	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\# Z > 2$
3	NAG	C	901	1	14,14,15	0.77	1 (7%)	17,19,21	0.72	0
3	NAG	B	901	1	14,14,15	0.54	0	17,19,21	0.46	0
3	NAG	D	901	1	14,14,15	0.93	1 (7%)	17,19,21	0.64	0
3	NAG	A	901	1	14,14,15	0.53	0	17,19,21	0.52	0

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	C	901	1	-	2/6/23/26	0/1/1/1
3	NAG	B	901	1	-	4/6/23/26	0/1/1/1
3	NAG	D	901	1	-	4/6/23/26	0/1/1/1
3	NAG	A	901	1	-	4/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	901	NAG	C1-C2	3.14	1.56	1.52
3	C	901	NAG	C1-C2	2.36	1.55	1.52

There are no bond angle outliers.

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	901	NAG	O5-C5-C6-O6
3	D	901	NAG	C4-C5-C6-O6
3	B	901	NAG	O5-C5-C6-O6
3	A	901	NAG	O5-C5-C6-O6
3	B	901	NAG	C4-C5-C6-O6
3	A	901	NAG	C4-C5-C6-O6
3	C	901	NAG	O5-C5-C6-O6
3	C	901	NAG	C3-C2-N2-C7
3	B	901	NAG	C1-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
3	D	901	NAG	C1-C2-N2-C7
3	A	901	NAG	C3-C2-N2-C7
3	B	901	NAG	C3-C2-N2-C7
3	A	901	NAG	C1-C2-N2-C7
3	D	901	NAG	C3-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	321/333 (96%)	-0.88	1 (0%) 90 75	52, 119, 208, 276	0
1	B	321/333 (96%)	-1.03	0 100 100	61, 111, 175, 250	0
1	C	321/333 (96%)	-0.90	0 100 100	50, 134, 211, 263	0
1	D	321/333 (96%)	-1.01	0 100 100	47, 110, 184, 261	0
2	E	247/261 (94%)	-1.03	0 100 100	59, 102, 162, 219	0
2	F	247/261 (94%)	-0.96	0 100 100	67, 116, 178, 307	0
2	G	247/261 (94%)	-0.97	0 100 100	56, 112, 177, 224	0
2	H	247/261 (94%)	-0.96	0 100 100	67, 116, 183, 262	0
All	All	2272/2376 (95%)	-0.97	1 (0%) 100 100	47, 114, 189, 307	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	288	GLN	2.3

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 oligosaccharides 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. 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(Å ²)	Q<0.9
3	NAG	A	901	14/15	0.97	0.07	135,141,145,146	0
3	NAG	B	901	14/15	0.97	0.05	103,120,123,125	0
3	NAG	C	901	14/15	0.98	0.04	88,103,107,107	0
4	CA	H	401	1/1	0.98	0.04	188,188,188,188	0
4	CA	E	401	1/1	0.99	0.04	133,133,133,133	0
4	CA	F	401	1/1	0.99	0.03	124,124,124,124	0
4	CA	G	401	1/1	0.99	0.04	98,98,98,98	0
3	NAG	D	901	14/15	0.99	0.03	85,101,110,110	0

6.5 Other polymers [i](#)

There are no such residues in this entry.