



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 05:04 AM UTC

PDB ID : 5JTW / pdb_00005jtw
Title : Crystal structure of complement C4b re-refined using iMDFF
Authors : Croll, T.I.; Andersen, G.R.
Deposited on : 2016-05-09
Resolution : 3.50 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

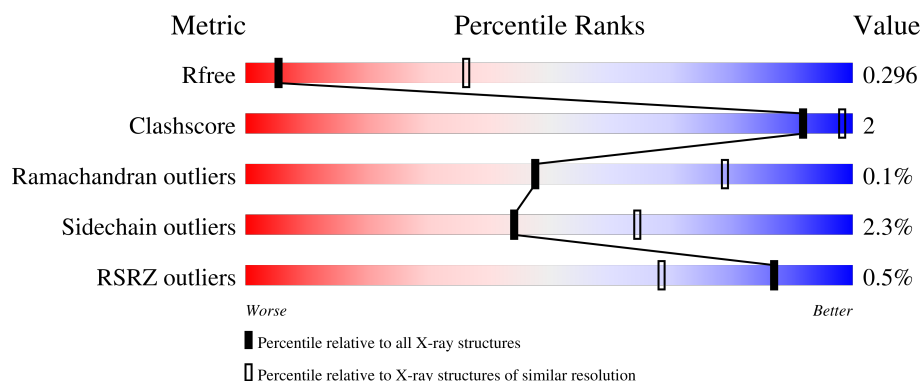
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	656	91%8% ..
1	D	656	94%5% .
2	B	690	% 85%6%8%
2	E	690	% 87%.8%
3	C	291	93%. .

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Mol	Chain	Length	Quality of chain
3	F	291	 90%6% •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 24312 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 Complement C4-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	651	Total	C	N	O	S	0	0	0
			5012	3185	872	939	16			
1	D	651	Total	C	N	O	S	0	0	0
			5012	3185	872	939	16			

- Molecule 2 is a protein called Complement C4-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	632	Total	C	N	O	S	0	0	0
			4876	3085	844	932	15			
2	E	632	Total	C	N	O	S	0	0	0
			4876	3085	844	932	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1201	SER	THR	conflict	UNP P0C0L4
E	1201	SER	THR	conflict	UNP P0C0L4

- Molecule 3 is a protein called Complement C4-A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	281	Total	C	N	O	S	0	0	0
			2240	1407	400	416	17			
3	F	281	Total	C	N	O	S	0	0	0
			2240	1407	400	416	17			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



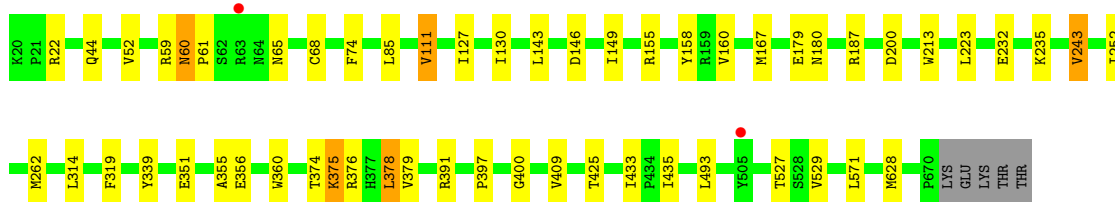
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		
4	E	1	Total	C	N	O	0	0
			14	8	1	5		

3 Residue-property plots

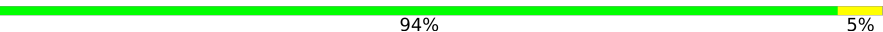
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

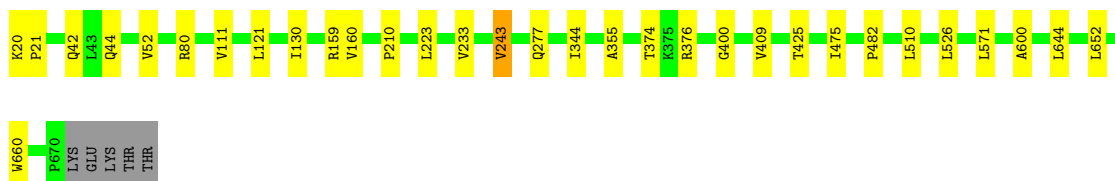
- Molecule 1: Complement C4-A

Chain A: 



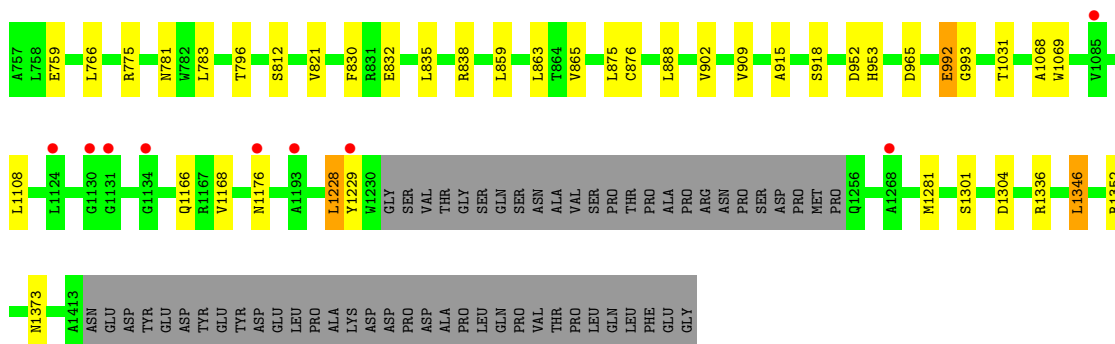
- Molecule 1: Complement C4-A

Chain D: 

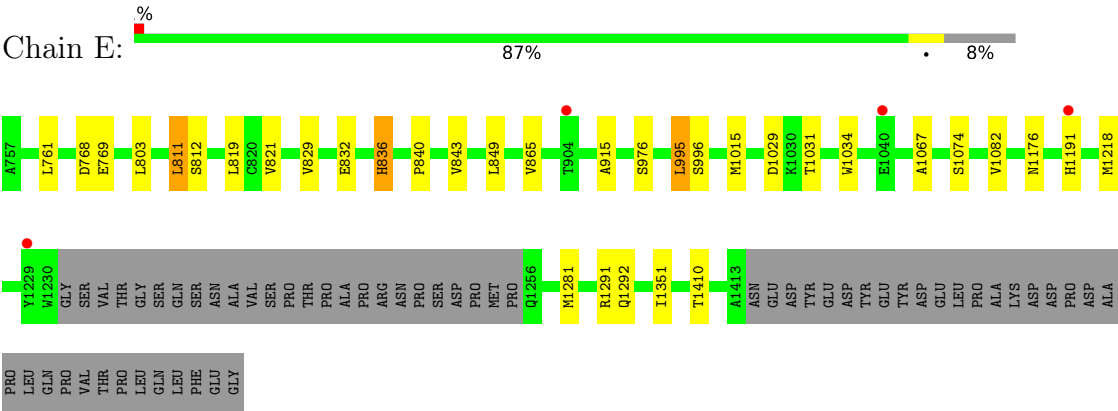


- Molecule 2: Complement C4-A

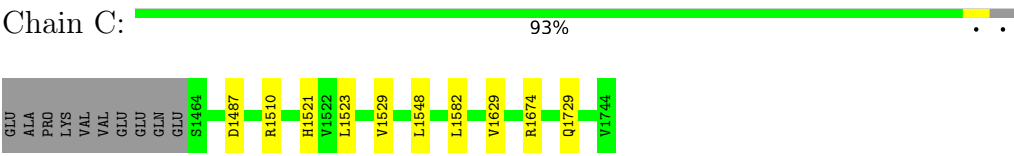
Chain B: 



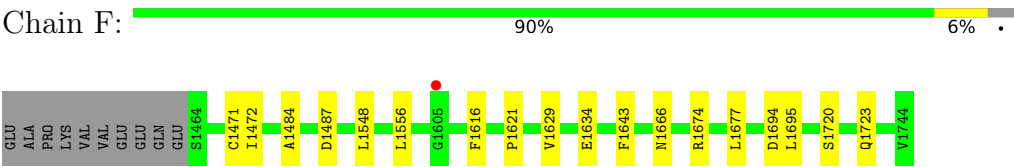
- Molecule 2: Complement C4-A



• Molecule 3: Complement C4-A



• Molecule 3: Complement C4-A



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	121.50Å 161.08Å 131.60Å 90.00° 107.26° 90.00°	Depositor
Resolution (Å)	49.38 – 3.50 49.38 – 3.50	Depositor EDS
% Data completeness (in resolution range)	98.7 (49.38-3.50) 98.6 (49.38-3.50)	Depositor EDS
R_{merge}	0.49	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 3.48Å)	Xtriage
Refinement program	PHENIX dev_2376	Depositor
R, R_{free}	0.248 , 0.293 0.248 , 0.296	Depositor DCC
R_{free} test set	2014 reflections (3.34%)	wwPDB-VP
Wilson B-factor (Å ²)	101.3	Xtriage
Anisotropy	0.345	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 98.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	24312	wwPDB-VP
Average B, all atoms (Å ²)	122.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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: 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	0.19	1/5128 (0.0%)	0.49	0/6961
1	D	0.15	0/5128	0.48	0/6961
2	B	0.16	0/4975	0.57	0/6765
2	E	0.15	0/4975	0.53	0/6765
3	C	0.16	0/2287	0.56	0/3090
3	F	0.15	0/2287	0.52	0/3090
All	All	0.16	1/24780 (0.0%)	0.52	0/33632

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	433	ILE	C-N	8.42	1.53	1.33

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	60	ASN	Peptide

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	5012	0	5025	24	1
1	D	5012	0	5025	11	0
2	B	4876	0	4845	20	1
2	E	4876	0	4845	17	0
3	C	2240	0	2184	2	0
3	F	2240	0	2184	6	2
4	A	14	0	13	0	0
4	B	14	0	13	0	0
4	D	14	0	13	0	0
4	E	14	0	13	0	0
All	All	24312	0	24160	75	2

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 2.

All (75) 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:A:167:MET:HE2	1:A:628:MET:HE2	1.66	0.76
2:E:1029:ASP:OD1	2:E:1034:TRP:NE1	2.29	0.61
1:D:374:THR:O	1:D:376:ARG:NH1	2.36	0.59
2:B:775:ARG:NH2	2:B:835:LEU:O	2.37	0.58
1:A:391:ARG:HA	1:A:397:PRO:HA	1.88	0.54
1:A:243:VAL:HG13	1:A:355:ALA:HB2	1.88	0.54
1:A:187:ARG:NH2	1:A:200:ASP:OD2	2.37	0.52
1:D:277:GLN:H	1:D:277:GLN:CD	2.16	0.52
1:A:146:ASP:OD1	2:B:781:ASN:ND2	2.43	0.51
1:A:179:GLU:CD	2:B:1352:ARG:HH22	2.19	0.51
1:D:243:VAL:HG13	1:D:355:ALA:HB2	1.93	0.50
1:A:262:MET:HE1	1:A:319:PHE:CG	2.46	0.50
1:A:339:TYR:OH	1:A:356:GLU:OE2	2.23	0.49
1:A:378:LEU:HD22	1:A:379:VAL:H	1.76	0.49
2:E:1015:MET:HE3	2:E:1082:VAL:HG22	1.95	0.48
1:A:375:LYS:HD2	1:A:375:LYS:H	1.79	0.48
2:B:992:GLU:CD	2:B:993:GLY:H	2.21	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:400:GLY:N	1:D:425:THR:OG1	2.39	0.47
1:A:374:THR:O	1:A:376:ARG:NH1	2.35	0.47
1:A:400:GLY:N	1:A:425:THR:OG1	2.45	0.47
2:E:843:VAL:HG21	2:E:849:LEU:HD22	1.96	0.47
2:E:1067:ALA:HB2	2:E:1074:SER:HA	1.96	0.47
2:E:1291:ARG:NH1	2:E:1292:GLN:OE1	2.43	0.47
1:A:59:ARG:CZ	1:A:65:ASN:HD21	2.28	0.47
1:A:111:VAL:HG13	1:A:127:ILE:HG23	1.98	0.46
2:B:876:CYS:HB3	2:B:902:VAL:HG13	1.98	0.46
2:B:859:LEU:HD12	2:B:863:LEU:HD11	1.97	0.46
2:B:766:LEU:HD22	2:B:918:SER:HA	1.97	0.45
2:B:1346:LEU:H	2:B:1346:LEU:HD23	1.80	0.45
2:B:1228:LEU:HD22	2:B:1228:LEU:HA	1.83	0.45
3:F:1471:CYS:C	3:F:1472:ILE:HD12	2.42	0.45
3:F:1720:SER:O	3:F:1723:GLN:NE2	2.49	0.45
2:B:830:PHE:CZ	2:B:832:GLU:HA	2.52	0.44
2:E:811:LEU:HD11	2:E:819:LEU:HG	1.99	0.44
3:F:1484:ALA:HB1	3:F:1556:LEU:HD21	1.98	0.44
1:D:210:PRO:HA	1:D:233:VAL:HG13	1.99	0.44
2:E:761:LEU:HD12	2:E:1410:THR:HG21	1.99	0.44
2:B:1301:SER:OG	2:B:1304:ASP:OD1	2.29	0.44
1:A:61:PRO:HD3	1:A:68:CYS:SG	2.57	0.44
2:E:811:LEU:HD13	2:E:821:VAL:HB	1.98	0.44
2:E:865:VAL:HG11	2:E:915:ALA:HB1	2.00	0.44
1:A:22:ARG:HB2	1:A:44:GLN:HB2	2.01	0.43
1:A:180:ASN:HA	1:A:213:TRP:CD2	2.54	0.43
1:A:351:GLU:OE1	2:B:838:ARG:NH2	2.49	0.43
1:D:42:GLN:NE2	1:D:44:GLN:OE1	2.35	0.43
1:A:223:LEU:H	1:A:223:LEU:HD23	1.84	0.43
2:B:992:GLU:CD	2:B:992:GLU:H	2.26	0.43
1:D:644:LEU:H	1:D:644:LEU:HD22	1.83	0.43
2:E:803:LEU:HA	2:E:829:VAL:HG23	2.01	0.43
2:B:1228:LEU:HD13	2:B:1229:TYR:H	1.84	0.43
2:E:840:PRO:HG3	2:E:849:LEU:HD21	2.00	0.42
1:A:262:MET:HB3	1:A:360:TRP:CH2	2.55	0.42
1:D:159:ARG:CZ	1:D:600:ALA:HB1	2.50	0.42
1:A:158:TYR:CE1	1:A:200:ASP:HB3	2.55	0.42
2:B:875:LEU:HG	2:B:909:VAL:HG11	2.01	0.42
2:B:865:VAL:HG11	2:B:915:ALA:HB1	2.01	0.42
2:E:768:ASP:O	2:E:769:GLU:HB3	2.20	0.42
3:F:1666:ASN:ND2	3:F:1694:ASP:OD2	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:832:GLU:H	2:E:832:GLU:CD	2.27	0.42
3:F:1643:PHE:CZ	3:F:1677:LEU:HB2	2.55	0.41
2:B:952:ASP:CG	2:B:953:HIS:H	2.28	0.41
3:C:1510:ARG:H	3:C:1510:ARG:HG2	1.66	0.41
2:B:1068:ALA:HB3	2:B:1069:TRP:CE3	2.54	0.41
2:B:1108:LEU:HD11	2:B:1168:VAL:HG22	2.01	0.41
1:A:149:ILE:HA	1:A:232:GLU:O	2.21	0.41
1:A:571:LEU:HD22	2:B:812:SER:HB2	2.03	0.41
3:F:1616:PHE:CE1	3:F:1621:PRO:HD2	2.55	0.41
1:D:571:LEU:HD22	2:E:812:SER:HB2	2.02	0.41
1:A:74:PHE:HB3	1:A:85:LEU:HD11	2.03	0.40
1:D:20:LYS:HB2	1:D:21:PRO:CD	2.51	0.40
3:C:1521:HIS:CE1	3:C:1523:LEU:HG	2.55	0.40
2:E:1218:MET:HE1	2:E:1281:MET:SD	2.61	0.40
2:E:995:LEU:HD22	2:E:996:SER:H	1.86	0.40
2:E:836:HIS:CD2	2:E:836:HIS:C	2.97	0.40
1:D:510:LEU:HD11	1:D:652:LEU:HG	2.03	0.40

All (2) 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)
2:B:1336:ARG:NH2	3:F:1695:LEU:O[2_555]	2.08	0.12
1:A:155:ARG:NH1	3:F:1634:GLU:OE2[2_555]	2.11	0.09

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	649/656 (99%)	623 (96%)	25 (4%)	1 (0%)	43 74
1	D	649/656 (99%)	626 (96%)	23 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	628/690 (91%)	579 (92%)	48 (8%)	1 (0%)	43	74
2	E	628/690 (91%)	598 (95%)	30 (5%)	0	100	100
3	C	279/291 (96%)	260 (93%)	19 (7%)	0	100	100
3	F	279/291 (96%)	260 (93%)	19 (7%)	0	100	100
All	All	3112/3274 (95%)	2946 (95%)	164 (5%)	2 (0%)	48	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
2	B	965	ASP

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	557/562 (99%)	541 (97%)	16 (3%)	37	60
1	D	557/562 (99%)	543 (98%)	14 (2%)	42	63
2	B	524/575 (91%)	511 (98%)	13 (2%)	42	63
2	E	524/575 (91%)	516 (98%)	8 (2%)	57	71
3	C	240/249 (96%)	233 (97%)	7 (3%)	37	60
3	F	240/249 (96%)	236 (98%)	4 (2%)	53	70
All	All	2642/2772 (95%)	2580 (98%)	62 (2%)	44	64

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	VAL
1	A	111	VAL
1	A	130	ILE
1	A	143	LEU
1	A	160	VAL

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Mol	Chain	Res	Type
1	A	235	LYS
1	A	243	VAL
1	A	252	ILE
1	A	314	LEU
1	A	375	LYS
1	A	378	LEU
1	A	409	VAL
1	A	435	ILE
1	A	493	LEU
1	A	527	THR
1	A	529	VAL
2	B	759	GLU
2	B	783	LEU
2	B	796	THR
2	B	821	VAL
2	B	888	LEU
2	B	992	GLU
2	B	1031	THR
2	B	1166	GLN
2	B	1176	ASN
2	B	1228	LEU
2	B	1281	MET
2	B	1346	LEU
2	B	1373	ASN
3	C	1487	ASP
3	C	1529	VAL
3	C	1548	LEU
3	C	1582	LEU
3	C	1629	VAL
3	C	1674	ARG
3	C	1729	GLN
1	D	52	VAL
1	D	80	ARG
1	D	111	VAL
1	D	121	LEU
1	D	130	ILE
1	D	160	VAL
1	D	223	LEU
1	D	243	VAL
1	D	344	ILE
1	D	409	VAL
1	D	475	ILE

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Mol	Chain	Res	Type
1	D	482	PRO
1	D	526	LEU
1	D	660	TRP
2	E	811	LEU
2	E	836	HIS
2	E	976	SER
2	E	995	LEU
2	E	1031	THR
2	E	1176	ASN
2	E	1191	HIS
2	E	1351	ILE
3	F	1487	ASP
3	F	1548	LEU
3	F	1629	VAL
3	F	1674	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	ASN
1	A	182	HIS
1	A	490	ASN
1	A	504	HIS
2	B	781	ASN
2	B	834	HIS
2	B	868	HIS
2	B	940	HIS
2	B	1112	GLN
2	B	1118	GLN
2	B	1166	GLN
2	B	1271	HIS
2	B	1275	HIS
2	B	1391	ASN
1	D	100	HIS
1	D	311	HIS
1	D	549	HIS
1	D	616	HIS
1	D	648	GLN
2	E	809	HIS
2	E	1176	ASN
2	E	1191	HIS
2	E	1342	HIS

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Mol	Chain	Res	Type
3	F	1467	HIS
3	F	1514	HIS
3	F	1731	ASN
3	F	1735	GLN

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](#)

4 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	$\# Z > 2$
4	NAG	A	1801	1	14,14,15	0.39	0	17,19,21	0.46	0
4	NAG	D	1801	1	14,14,15	0.36	0	17,19,21	0.42	0
4	NAG	E	1501	2	14,14,15	0.44	0	17,19,21	0.45	0
4	NAG	B	1501	2	14,14,15	0.53	0	17,19,21	0.49	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
4	NAG	A	1801	1	-	0/6/23/26	0/1/1/1
4	NAG	D	1801	1	-	1/6/23/26	0/1/1/1
4	NAG	E	1501	2	-	0/6/23/26	0/1/1/1
4	NAG	B	1501	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	1801	NAG	C1-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

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	651/656 (99%)	-0.12	2 (0%) 90 76	42, 108, 160, 194	0
1	D	651/656 (99%)	-0.16	0 100 100	51, 93, 137, 177	0
2	B	632/690 (91%)	0.15	9 (1%) 73 48	42, 156, 239, 275	0
2	E	632/690 (91%)	0.05	4 (0%) 85 65	77, 124, 180, 246	0
3	C	281/291 (96%)	0.10	0 100 100	59, 120, 190, 225	0
3	F	281/291 (96%)	0.04	1 (0%) 88 72	93, 130, 186, 219	0
All	All	3128/3274 (95%)	-0.00	16 (0%) 87 68	42, 115, 200, 275	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	505	TYR	4.1
2	E	1229	TYR	3.4
2	B	1268	ALA	3.2
2	E	1191	HIS	2.9
2	B	1131	GLY	2.9
3	F	1605	GLY	2.7
2	B	1229	TYR	2.6
1	A	63	ARG	2.5
2	B	1193	ALA	2.4
2	E	904	THR	2.3
2	B	1130	GLY	2.2
2	B	1176	ASN	2.2
2	E	1040	GLU	2.1
2	B	1124	LEU	2.0
2	B	1085	VAL	2.0
2	B	1134	GLY	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 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($\text{\AA}^2$)	Q<0.9
4	NAG	A	1801	14/15	0.56	0.12	90,148,184,193	0
4	NAG	D	1801	14/15	0.58	0.10	132,158,172,184	0
4	NAG	E	1501	14/15	0.69	0.09	142,159,173,178	0
4	NAG	B	1501	14/15	0.75	0.08	124,143,158,158	0

6.5 Other polymers [i](#)

There are no such residues in this entry.