



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:42 AM UTC

PDB ID : 8BCG / pdb_00008bcg
Title : Human Brr2 Helicase Region in complex with C-tail deleted Jab1 and compound 86
Authors : Vester, K.; Loll, B.; Wahl, M.C.
Deposited on : 2022-10-15
Resolution : 2.39 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

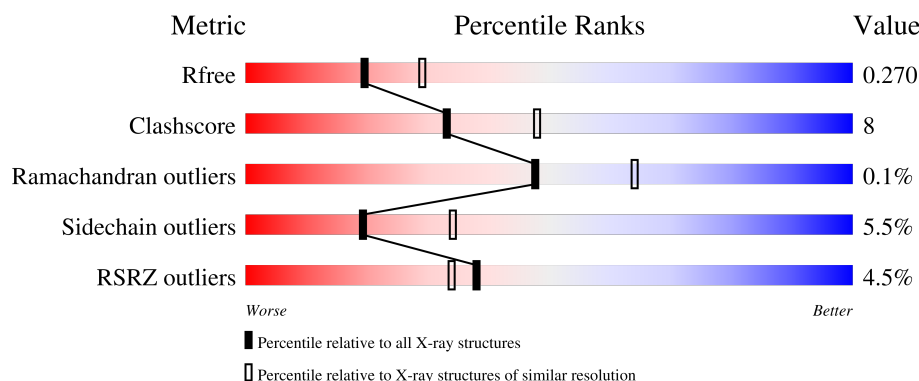
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	B	1747	
2	J	263	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16360 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 U5 small nuclear ribonucleoprotein 200 kDa helicase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	1725	Total	C	N	O	S	0	0	0
			13870	8865	2373	2560	72			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	390	GLY	-	expression tag	UNP O75643
B	391	ALA	-	expression tag	UNP O75643
B	392	GLU	-	expression tag	UNP O75643
B	393	PHE	-	expression tag	UNP O75643

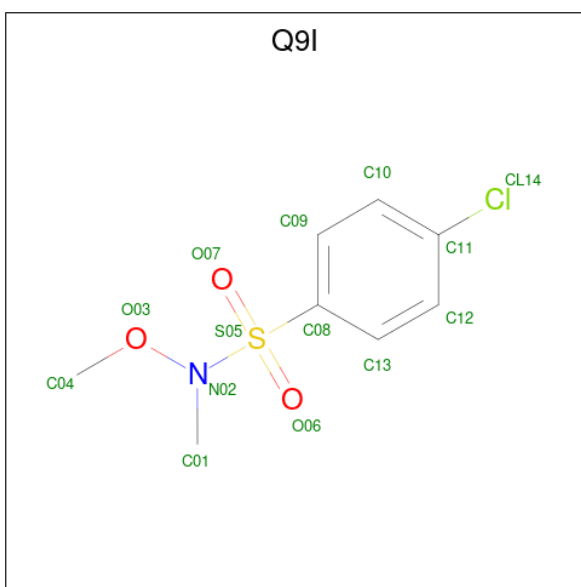
- Molecule 2 is a protein called Pre-mRNA-processing-splicing factor 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	J	260	Total	C	N	O	S	0	0	0
			2104	1345	362	385	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
J	2058	GLY	-	expression tag	UNP Q6P2Q9
J	2059	PRO	-	expression tag	UNP Q6P2Q9
J	2060	LEU	-	expression tag	UNP Q6P2Q9
J	2061	GLY	-	expression tag	UNP Q6P2Q9
J	2062	SER	-	expression tag	UNP Q6P2Q9
J	2063	MET	-	expression tag	UNP Q6P2Q9

- Molecule 3 is 4-chloranyl- {N}-methoxy- {N}-methyl-benzenesulfonamide (CCD ID: Q9I) (formula: C₈H₁₀ClNO₃S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	Cl	N	O	S	0	0
			14	8	1	1	3	1		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	C O	0	0
			4	2 2		
4	B	1	Total	C O	0	0
			4	2 2		
4	B	1	Total	C O	0	0
			4	2 2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0

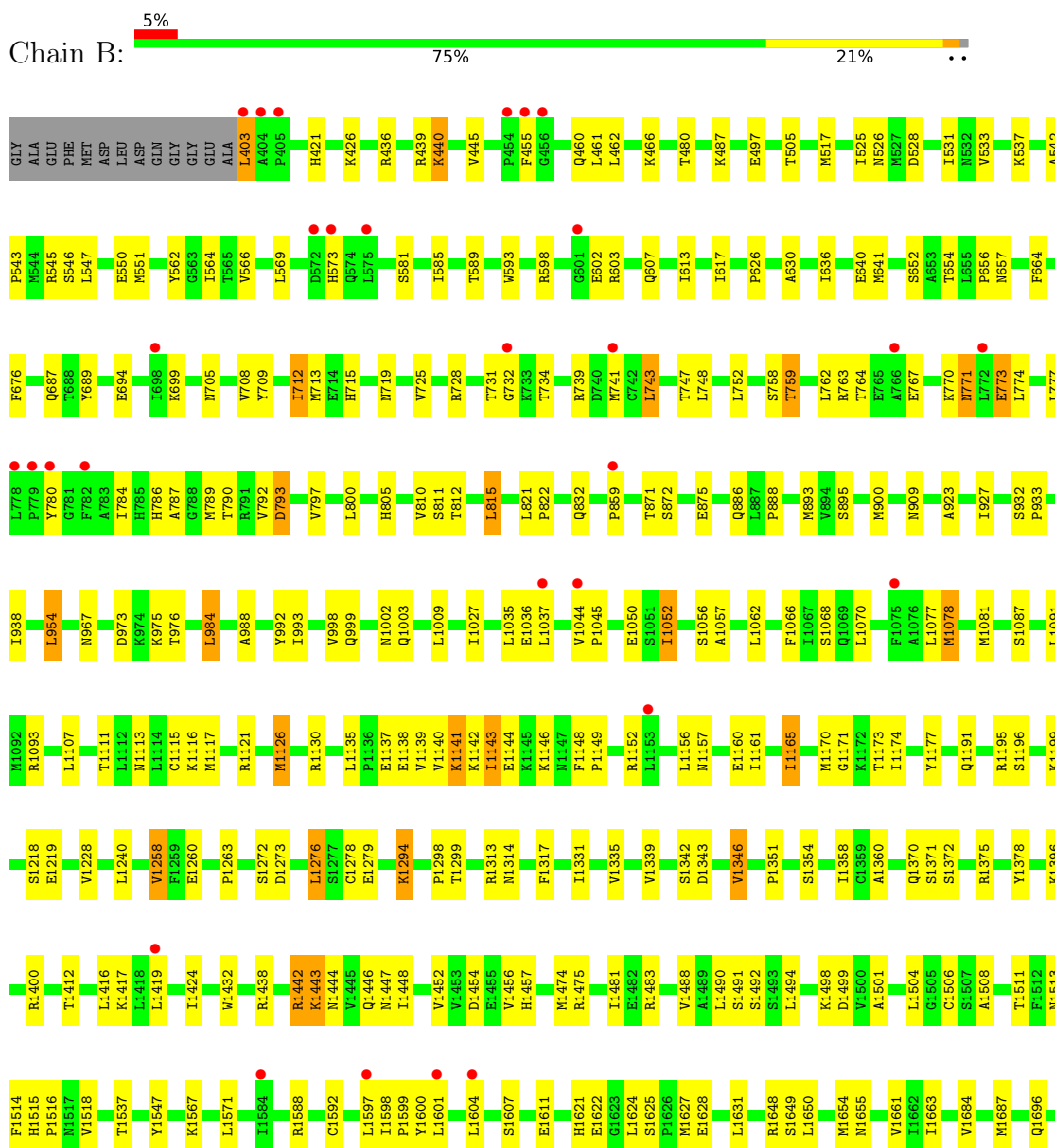
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	287	Total O 287 287	0	0
5	J	49	Total O 49 49	0	0

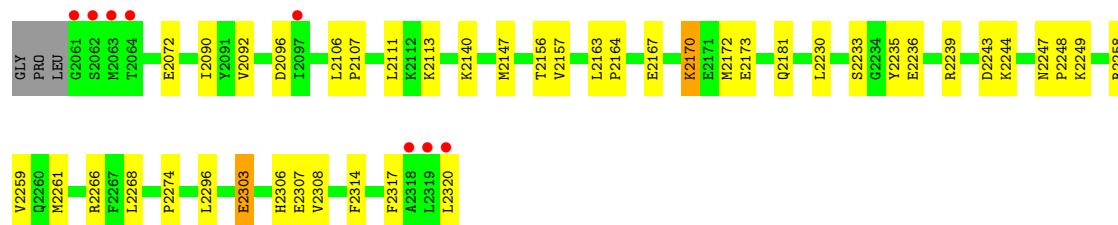
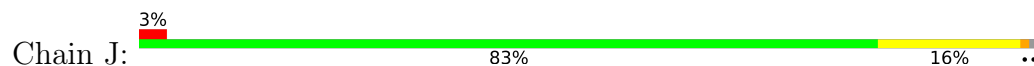
3 Residue-property plots

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: U5 small nuclear ribonucleoprotein 200 kDa helicase



- Molecule 2: Pre-mRNA-processing-splicing factor 8



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.92Å 118.92Å 188.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.06 – 2.39 46.06 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.06-2.39) 99.8 (46.06-2.39)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.214 , 0.269 0.214 , 0.270	Depositor DCC
R_{free} test set	2101 reflections (2.35%)	wwPDB-VP
Wilson B-factor (Å ²)	48.3	Xtriage
Anisotropy	0.183	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16360	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% 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: EDO, Q9I

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	B	0.33	0/14164	0.51	0/19191
2	J	0.33	0/2170	0.50	0/2953
All	All	0.33	0/16334	0.51	0/22144

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	B	13870	0	14018	232	0
2	J	2104	0	2042	19	0
3	B	14	0	0	1	0
4	B	36	0	54	1	0
5	B	287	0	0	7	0
5	J	49	0	0	1	0
All	All	16360	0	16114	250	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 8.

All (250) 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:993:ILE:HD12	1:B:1091:LEU:HD23	1.51	0.90
1:B:1879:LEU:HD11	1:B:1882:PRO:HB3	1.58	0.85
1:B:440:LYS:H	1:B:440:LYS:HD3	1.45	0.82
1:B:1747:ASN:HB2	1:B:1808:MET:HE3	1.62	0.81
1:B:421:HIS:NE2	1:B:875:GLU:OE1	2.16	0.74
1:B:2006:ASP:HA	1:B:2009:ARG:HH11	1.53	0.73
1:B:1836:LEU:HD22	1:B:1930:LEU:HD21	1.71	0.72
1:B:728:ARG:HE	1:B:787:ALA:HB3	1.54	0.72
1:B:762:LEU:HD21	1:B:805:HIS:HB3	1.71	0.71
1:B:1475:ARG:HD2	1:B:1504:LEU:HA	1.72	0.70
1:B:2043:ARG:NH2	1:B:2084:LEU:O	2.26	0.69
1:B:763:ARG:NH2	1:B:764:THR:OG1	2.26	0.68
1:B:1351:PRO:HG3	1:B:1516:PRO:HA	1.77	0.66
1:B:984:LEU:HD21	1:B:1002:ASN:HB2	1.77	0.64
1:B:1732:MET:HE3	1:B:1788:LEU:HG	1.78	0.64
1:B:1696:GLN:OE1	5:B:2301:HOH:O	2.14	0.63
1:B:790:THR:HG22	1:B:792:VAL:H	1.64	0.62
1:B:1604:LEU:HD23	1:B:1628:GLU:HG2	1.82	0.62
1:B:1684:VAL:HA	1:B:1687:MET:HE2	1.81	0.62
1:B:1138:GLU:H	1:B:1138:GLU:CD	2.07	0.61
1:B:1066:PHE:HD1	1:B:1081:MET:HE3	1.66	0.60
1:B:1456:VAL:HG12	1:B:1491:SER:HB2	1.83	0.60
1:B:436:ARG:HG3	1:B:445:VAL:HG22	1.83	0.60
2:J:2235:TYR:O	2:J:2239:ARG:HG2	2.01	0.60
1:B:1906:ALA:O	1:B:1910:SER:OG	2.20	0.60
1:B:1804:ILE:HG12	1:B:1810:VAL:HG12	1.84	0.59
1:B:2006:ASP:OD1	1:B:2009:ARG:NH1	2.34	0.59
1:B:497:GLU:HG3	5:B:2330:HOH:O	2.01	0.59
1:B:1663:ILE:HD12	1:B:1704:ILE:HG12	1.84	0.59
1:B:793:ASP:OD1	1:B:793:ASP:N	2.36	0.59
1:B:1499:ASP:OD2	1:B:1763:ARG:NH1	2.32	0.59
1:B:1299:THR:HG22	5:B:2583:HOH:O	2.03	0.58
1:B:2051:VAL:HG13	1:B:2113:TYR:CZ	2.38	0.58
1:B:1452:VAL:HG22	1:B:1488:VAL:HB	1.85	0.58
1:B:1764:MET:HE3	1:B:1773:LEU:HD11	1.85	0.57
1:B:975:LYS:HG3	1:B:976:THR:HG23	1.87	0.57
1:B:1442:ARG:HE	1:B:1442:ARG:HA	1.70	0.56
1:B:1142:LYS:O	1:B:1146:LYS:HG2	2.06	0.56
1:B:1314:ASN:HB3	1:B:1317:PHE:HB2	1.87	0.56
1:B:1066:PHE:CZ	1:B:1121:ARG:HD3	2.39	0.56
1:B:1515:HIS:HB3	1:B:1516:PRO:HD2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:598:ARG:HH22	4:B:2206:EDO:H11	1.69	0.56
1:B:1592:CYS:SG	1:B:1597:LEU:HD13	2.47	0.55
1:B:1514:PHE:HB3	1:B:1518:VAL:HG21	1.87	0.55
1:B:1444:ASN:HA	1:B:1447:ASN:ND2	2.21	0.55
1:B:1967:THR:HG22	1:B:1970:HIS:ND1	2.22	0.55
1:B:1814:ASN:HA	1:B:1817:MET:HE3	1.88	0.55
1:B:1598:ILE:HA	1:B:1601:LEU:HB2	1.88	0.55
1:B:1191:GLN:OE1	1:B:1199:LYS:HE3	2.08	0.54
1:B:1346:VAL:HG13	1:B:1488:VAL:HG13	1.90	0.54
1:B:1739:GLU:HA	1:B:1742:THR:HG22	1.89	0.54
1:B:1648:ARG:HG3	1:B:1649:SER:N	2.22	0.54
1:B:1843:ARG:HD3	1:B:1877:HIS:ND1	2.22	0.54
1:B:543:PRO:HD2	1:B:547:LEU:HD23	1.90	0.53
1:B:664:PHE:HB2	1:B:927:ILE:HD13	1.91	0.53
1:B:1432:TRP:HE1	1:B:1474:MET:HE3	1.73	0.53
1:B:1161:ILE:HD11	1:B:1171:GLY:HA2	1.90	0.53
1:B:1858:ILE:HB	1:B:1887:PRO:HB3	1.90	0.53
1:B:1864:GLU:HB3	1:B:1868:LEU:HD23	1.90	0.53
2:J:2236:GLU:OE2	2:J:2239:ARG:NH1	2.42	0.53
2:J:2106:LEU:HD12	2:J:2107:PRO:HD2	1.91	0.53
1:B:1416:LEU:HD22	1:B:1442:ARG:HD3	1.90	0.52
1:B:2015:PRO:HG2	1:B:2116:CYS:SG	2.49	0.52
1:B:2037:VAL:HG11	1:B:2068:ILE:HD12	1.90	0.52
1:B:593:TRP:HE1	3:B:2201:Q9I:C01	2.22	0.52
1:B:1107:LEU:O	1:B:1111:THR:HG23	2.09	0.52
1:B:1866:ASN:HA	1:B:1869:ARG:HB2	1.92	0.52
1:B:725:VAL:HG21	1:B:731:THR:HA	1.92	0.52
1:B:1037:LEU:HB3	1:B:1052:ILE:HD11	1.92	0.52
1:B:2074:ASN:ND2	1:B:2074:ASN:O	2.43	0.52
1:B:1343:ASP:OD1	1:B:1370:GLN:NE2	2.42	0.52
1:B:1973:ARG:HG2	1:B:1996:LEU:HD11	1.92	0.52
1:B:2066:VAL:O	1:B:2079:ILE:HA	2.10	0.52
1:B:815:LEU:HD21	1:B:821:LEU:HD22	1.92	0.51
1:B:1851:ASN:HA	1:B:1888:HIS:ND1	2.26	0.51
1:B:1298:PRO:HB3	1:B:1515:HIS:CG	2.46	0.51
1:B:1156:LEU:HD12	1:B:1161:ILE:HG22	1.92	0.50
1:B:694:GLU:HG2	1:B:699:LYS:HG3	1.94	0.50
1:B:1149:PRO:O	1:B:1152:ARG:HG2	2.11	0.50
1:B:1360:ALA:HB2	1:B:1490:LEU:HD11	1.93	0.50
1:B:654:THR:HG21	1:B:676:PHE:O	2.12	0.50
1:B:2073:SER:HB2	1:B:2075:SER:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2051:VAL:HG11	1:B:2112:ALA:HB1	1.93	0.49
1:B:439:ARG:HH21	1:B:440:LYS:HZ3	1.60	0.49
1:B:630:ALA:HB2	1:B:900:MET:HE3	1.93	0.49
1:B:933:PRO:HB2	1:B:938:ILE:HB	1.94	0.49
1:B:1501:ALA:HB1	1:B:1506:CYS:HB2	1.95	0.49
1:B:1195:ARG:NH1	1:B:1260:GLU:OE2	2.46	0.49
1:B:709:TYR:O	1:B:712:ILE:HG22	2.13	0.49
1:B:786:HIS:CD2	1:B:789:MET:HE3	2.48	0.49
1:B:1943:MET:HG2	1:B:2065:TRP:CE3	2.48	0.49
1:B:2046:GLU:HG2	1:B:2086:GLN:HB2	1.95	0.49
1:B:1170:MET:HB2	1:B:1173:THR:HB	1.95	0.49
1:B:1078:MET:HA	1:B:1078:MET:HE3	1.94	0.48
1:B:1219:GLU:HG3	5:B:2319:HOH:O	2.12	0.48
1:B:1948:MET:HE3	1:B:1954:TRP:HA	1.94	0.48
2:J:2181:GLN:NE2	5:J:2402:HOH:O	2.46	0.48
1:B:1135:LEU:HD13	1:B:1177:TYR:CD2	2.49	0.48
1:B:1879:LEU:HD22	1:B:1893:LEU:HD21	1.94	0.48
1:B:1140:VAL:O	1:B:1144:GLU:HG2	2.14	0.48
1:B:1621:HIS:CE1	1:B:1624:LEU:HG	2.48	0.48
1:B:1481:ILE:O	1:B:1483:ARG:HG2	2.14	0.48
1:B:822:PRO:HG3	1:B:859:PRO:HD3	1.96	0.48
1:B:531:ILE:HD13	1:B:562:TYR:O	2.13	0.48
1:B:566:VAL:HG22	1:B:585:ILE:HB	1.95	0.48
1:B:1066:PHE:CE2	1:B:1121:ARG:HD3	2.49	0.47
1:B:1165:ILE:HD13	1:B:1174:ILE:HD13	1.95	0.47
1:B:705:ASN:HA	1:B:708:VAL:HG12	1.96	0.47
1:B:1062:LEU:HD21	1:B:1077:LEU:HD22	1.96	0.47
1:B:1137:GLU:HG2	1:B:1141:LYS:HD3	1.95	0.47
1:B:1904:LEU:HD13	1:B:1908:LEU:HD13	1.95	0.47
2:J:2092:VAL:HG13	2:J:2261:MET:HE3	1.96	0.47
1:B:617:ILE:HG22	1:B:652:SER:HB2	1.95	0.47
1:B:1126:MET:HB3	1:B:1130:ARG:HD2	1.96	0.47
1:B:641:MET:HE1	1:B:909:ASN:HA	1.97	0.47
1:B:505:THR:HG22	5:B:2388:HOH:O	2.14	0.47
1:B:1035:LEU:HD23	1:B:1036:GLU:H	1.80	0.47
1:B:487:LYS:HD2	1:B:676:PHE:HE1	1.80	0.47
1:B:1156:LEU:HB3	1:B:1160:GLU:HB2	1.97	0.47
1:B:1882:PRO:HB2	1:B:1884:PHE:CZ	2.50	0.47
1:B:2065:TRP:CZ3	1:B:2111:ASP:HB3	2.50	0.47
1:B:2109:MET:HE3	1:B:2109:MET:HB3	1.86	0.47
1:B:743:LEU:HD22	1:B:748:LEU:HD13	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1196:SER:HA	1:B:1258:VAL:HG12	1.97	0.47
1:B:871:THR:OG1	1:B:872:SER:N	2.47	0.46
1:B:728:ARG:O	1:B:731:THR:HG22	2.16	0.46
1:B:1156:LEU:HB2	1:B:1161:ILE:HG23	1.97	0.46
1:B:421:HIS:ND1	5:B:2306:HOH:O	2.36	0.46
1:B:767:GLU:O	1:B:770:LYS:HE3	2.14	0.46
1:B:800:LEU:O	1:B:805:HIS:HB2	2.14	0.46
1:B:1457:HIS:HB3	1:B:1494:LEU:HD21	1.97	0.46
1:B:752:LEU:HD13	1:B:759:THR:HG23	1.98	0.46
1:B:771:ASN:HB2	1:B:774:LEU:HB3	1.98	0.46
1:B:1943:MET:HE1	1:B:2067:VAL:HG21	1.98	0.46
1:B:758:SER:O	1:B:762:LEU:HG	2.16	0.46
1:B:1396:LYS:O	1:B:1400:ARG:HB2	2.15	0.46
1:B:1870:GLN:O	1:B:1873:GLN:HB3	2.16	0.46
2:J:2140:LYS:NZ	2:J:2173:GLU:OE2	2.34	0.46
1:B:1331:ILE:HD12	1:B:1354:SER:HB3	1.98	0.46
1:B:687:GLN:OE1	1:B:689:TYR:OH	2.12	0.45
2:J:2243:ASP:OD2	2:J:2248:PRO:HA	2.16	0.45
1:B:709:TYR:HB2	1:B:741:MET:HE3	1.97	0.45
1:B:1868:LEU:HD13	1:B:1871:LEU:HD12	1.98	0.45
1:B:1975:THR:HG22	1:B:1980:GLU:OE1	2.17	0.45
1:B:886:GLN:O	1:B:888:PRO:HD3	2.16	0.45
1:B:1218:SER:HB2	1:B:1240:LEU:HD11	1.99	0.45
1:B:1779:ARG:HB3	1:B:1779:ARG:NH1	2.31	0.45
1:B:2036:VAL:HG22	1:B:2093:ASP:HB3	1.99	0.45
1:B:1045:PRO:HD3	2:J:2317:PHE:CD2	2.51	0.45
1:B:1148:PHE:HD2	1:B:1152:ARG:HD3	1.81	0.45
1:B:1887:PRO:O	1:B:1891:THR:HG23	2.16	0.45
1:B:1378:TYR:OH	1:B:1454:ASP:OD2	2.23	0.45
1:B:636:ILE:O	1:B:640:GLU:HG2	2.17	0.44
1:B:542:ALA:HB1	1:B:547:LEU:HD23	1.99	0.44
1:B:1113:ASN:O	1:B:1117:MET:HG3	2.18	0.44
1:B:1622:GLU:OE1	1:B:1622:GLU:N	2.41	0.44
2:J:2090:ILE:HG21	2:J:2111:LEU:HD21	1.98	0.44
1:B:1513:ASN:ND2	5:B:2303:HOH:O	2.28	0.44
1:B:2039:VAL:HG21	1:B:2106:LEU:HD11	1.99	0.44
1:B:893:MET:HE3	1:B:900:MET:SD	2.58	0.44
1:B:999:GLN:OE1	1:B:1003:GLN:NE2	2.30	0.44
1:B:1442:ARG:HH12	1:B:1444:ASN:ND2	2.15	0.44
1:B:1607:SER:O	1:B:1611:GLU:HG3	2.18	0.44
1:B:2029:ILE:HG13	1:B:2124:VAL:CG1	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1068:SER:HB2	1:B:1070:LEU:HD13	1.99	0.44
1:B:2051:VAL:HG22	1:B:2062:GLU:HG2	2.00	0.44
1:B:1126:MET:HG2	1:B:1130:ARG:NH1	2.33	0.44
1:B:1846:ILE:O	1:B:1850:SER:OG	2.32	0.44
1:B:1878:LYS:HA	1:B:1878:LYS:HD3	1.53	0.44
1:B:1985:ILE:HG22	1:B:1993:ARG:HH21	1.82	0.44
2:J:2163:LEU:HD23	2:J:2163:LEU:HA	1.80	0.44
1:B:1272:SER:HB2	1:B:1278:CYS:SG	2.58	0.43
1:B:1442:ARG:HA	1:B:1442:ARG:NE	2.33	0.43
1:B:1116:LYS:HB2	1:B:1276:LEU:HD13	2.00	0.43
1:B:517:MET:HE3	1:B:613:ILE:HD12	2.00	0.43
1:B:550:GLU:OE1	1:B:551:MET:HE2	2.18	0.43
1:B:1062:LEU:HD22	1:B:1081:MET:HB2	2.00	0.43
1:B:1157:ASN:ND2	1:B:1160:GLU:OE2	2.49	0.43
1:B:1442:ARG:HH21	1:B:1443:LYS:HE2	1.82	0.43
2:J:2167:GLU:HA	2:J:2170:LYS:HE3	2.00	0.43
1:B:1093:ARG:HD2	1:B:1115:CYS:SG	2.57	0.43
2:J:2096:ASP:OD1	2:J:2096:ASP:N	2.52	0.43
1:B:739:ARG:HH11	1:B:780:TYR:HE2	1.66	0.43
1:B:1142:LYS:HE2	1:B:1142:LYS:HB2	1.92	0.43
2:J:2303:GLU:CD	2:J:2303:GLU:H	2.26	0.43
2:J:2320:LEU:HD23	2:J:2320:LEU:HA	1.80	0.43
1:B:562:TYR:HB2	1:B:564:ILE:HD12	2.00	0.43
1:B:1567:LYS:O	1:B:1571:LEU:HG	2.19	0.43
1:B:1598:ILE:N	1:B:1599:PRO:HD2	2.33	0.43
1:B:771:ASN:HB2	1:B:774:LEU:CB	2.49	0.43
2:J:2113:LYS:HE2	2:J:2113:LYS:HB2	1.86	0.43
1:B:602:GLU:HG3	1:B:603:ARG:H	1.84	0.43
1:B:975:LYS:HG3	1:B:976:THR:N	2.34	0.43
1:B:1335:VAL:O	1:B:1339:VAL:HG23	2.19	0.43
1:B:1419:LEU:HG	1:B:1444:ASN:HB3	2.00	0.43
1:B:531:ILE:O	1:B:533:VAL:HG13	2.19	0.42
1:B:784:ILE:HG13	1:B:789:MET:HE1	2.01	0.42
1:B:1191:GLN:HB3	1:B:1770:TYR:CD1	2.54	0.42
1:B:1732:MET:HE2	1:B:1732:MET:HB2	1.96	0.42
1:B:1872:ALA:HB1	1:B:1879:LEU:HD21	2.00	0.42
1:B:967:ASN:HB3	1:B:999:GLN:HB2	2.01	0.42
1:B:1627:MET:HE3	1:B:1627:MET:HB3	1.84	0.42
1:B:973:ASP:OD1	1:B:975:LYS:HG2	2.20	0.42
1:B:1139:VAL:O	1:B:1143:ILE:HD12	2.19	0.42
1:B:1228:VAL:CG2	1:B:1263:PRO:HB2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1375:ARG:HB3	1:B:1448:ILE:HD12	2.02	0.42
1:B:1650:LEU:O	1:B:1654:MET:HE3	2.19	0.42
1:B:1661:VAL:O	1:B:1702:CYS:HA	2.19	0.42
1:B:793:ASP:O	1:B:797:VAL:HG23	2.20	0.42
1:B:923:ALA:O	1:B:927:ILE:HD12	2.20	0.42
1:B:1891:THR:HG22	1:B:1915:ILE:HD11	2.01	0.42
1:B:626:PRO:HG3	1:B:893:MET:HA	2.01	0.42
1:B:731:THR:HG23	1:B:732:GLY:N	2.35	0.42
1:B:988:ALA:HB2	1:B:998:VAL:HG11	2.02	0.42
1:B:1195:ARG:HD2	1:B:1294:LYS:HG3	2.00	0.42
1:B:1498:LYS:HD2	1:B:1498:LYS:HA	1.81	0.42
1:B:2037:VAL:HG11	1:B:2068:ILE:CD1	2.49	0.42
1:B:439:ARG:HE	1:B:440:LYS:NZ	2.18	0.41
1:B:462:LEU:HD11	1:B:466:LYS:HB2	2.02	0.41
1:B:1148:PHE:CD2	1:B:1152:ARG:HD3	2.55	0.41
2:J:2307:GLU:HG3	2:J:2314:PHE:CE1	2.55	0.41
1:B:569:LEU:HB3	1:B:573:HIS:NE2	2.35	0.41
2:J:2164:PRO:HB3	2:J:2296:LEU:HD11	2.02	0.41
1:B:1967:THR:HG22	1:B:1970:HIS:CE1	2.55	0.41
1:B:709:TYR:CZ	1:B:713:MET:HE3	2.56	0.41
1:B:2029:ILE:HD12	1:B:2035:VAL:HG22	2.03	0.41
1:B:1050:GLU:HB2	1:B:1057:ALA:HB2	2.03	0.41
1:B:1438:ARG:O	1:B:1442:ARG:HB2	2.20	0.41
1:B:1858:ILE:HG13	1:B:1891:THR:CG2	2.50	0.41
1:B:715:HIS:HB3	1:B:719:ASN:HB2	2.01	0.41
1:B:784:ILE:HA	1:B:810:VAL:O	2.20	0.41
1:B:1417:LYS:HA	1:B:1417:LYS:HD3	1.77	0.41
1:B:1600:TYR:HD1	1:B:1631:LEU:HD22	1.86	0.41
2:J:2147:MET:O	2:J:2274:PRO:HD3	2.20	0.41
1:B:542:ALA:O	1:B:589:THR:HA	2.21	0.41
1:B:811:SER:OG	1:B:812:THR:N	2.54	0.41
1:B:1547:TYR:OH	1:B:1588:ARG:NH2	2.54	0.41
1:B:1899:LEU:HD22	1:B:1953:MET:O	2.21	0.41
1:B:1927:VAL:O	1:B:1931:SER:HB2	2.21	0.41
1:B:1945:LEU:O	1:B:1949:VAL:HG23	2.21	0.41
1:B:1093:ARG:NH2	1:B:1273:ASP:OD1	2.54	0.40
1:B:439:ARG:HE	1:B:440:LYS:HZ3	1.70	0.40
2:J:2306:HIS:ND1	2:J:2308:VAL:HG22	2.37	0.40
1:B:773:GLU:O	1:B:777:LEU:HD12	2.21	0.40
1:B:1733:HIS:HB3	1:B:1796:LEU:HD21	2.04	0.40
1:B:403:LEU:HB2	1:B:954:LEU:HD23	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1375:ARG:HB3	1:B:1448:ILE:CD1	2.51	0.40
1:B:656:PRO:O	1:B:657:ASN:HB2	2.22	0.40
1:B:1498:LYS:HE2	1:B:1508:ALA:HA	2.04	0.40
1:B:1781:LEU:O	1:B:1785:LEU:HG	2.21	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	B	1723/1747 (99%)	1648 (96%)	73 (4%)	2 (0%)	48	64
2	J	258/263 (98%)	245 (95%)	13 (5%)	0	100	100
All	All	1981/2010 (99%)	1893 (96%)	86 (4%)	2 (0%)	48	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2086	GLN
1	B	1881	ASN

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	1544/1560 (99%)	1461 (95%)	83 (5%)	20	35
2	J	234/236 (99%)	219 (94%)	15 (6%)	16	28
All	All	1778/1796 (99%)	1680 (94%)	98 (6%)	19	34

All (98) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	403	LEU
1	B	426	LYS
1	B	440	LYS
1	B	455	PHE
1	B	460	GLN
1	B	461	LEU
1	B	480	THR
1	B	525	ILE
1	B	526	ASN
1	B	528	ASP
1	B	537	LYS
1	B	545	ARG
1	B	546	SER
1	B	581	SER
1	B	607	GLN
1	B	712	ILE
1	B	734	THR
1	B	743	LEU
1	B	747	THR
1	B	759	THR
1	B	771	ASN
1	B	773	GLU
1	B	793	ASP
1	B	815	LEU
1	B	832	GLN
1	B	895	SER
1	B	932	SER
1	B	954	LEU
1	B	984	LEU
1	B	992	TYR
1	B	1009	LEU
1	B	1027	ILE

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Mol	Chain	Res	Type
1	B	1044	VAL
1	B	1052	ILE
1	B	1056	SER
1	B	1078	MET
1	B	1087	SER
1	B	1126	MET
1	B	1141	LYS
1	B	1143	ILE
1	B	1165	ILE
1	B	1258	VAL
1	B	1276	LEU
1	B	1279	GLU
1	B	1294	LYS
1	B	1313	ARG
1	B	1342	SER
1	B	1346	VAL
1	B	1358	ILE
1	B	1371	SER
1	B	1372	SER
1	B	1412	THR
1	B	1424	ILE
1	B	1442	ARG
1	B	1443	LYS
1	B	1446	GLN
1	B	1492	SER
1	B	1511	THR
1	B	1537	THR
1	B	1625	SER
1	B	1655	ASN
1	B	1744	THR
1	B	1754	TYR
1	B	1779	ARG
1	B	1781	LEU
1	B	1849	ILE
1	B	1861	ARG
1	B	1866	ASN
1	B	1871	LEU
1	B	1879	LEU
1	B	1910	SER
1	B	1931	SER
1	B	1953	MET
1	B	1958	SER

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Mol	Chain	Res	Type
1	B	1960	LEU
1	B	1966	PHE
1	B	1971	ILE
1	B	1974	CYS
1	B	1976	ASP
1	B	2030	ARG
1	B	2068	ILE
1	B	2090	VAL
1	B	2105	THR
2	J	2072	GLU
2	J	2156	THR
2	J	2157	VAL
2	J	2170	LYS
2	J	2172	MET
2	J	2230	LEU
2	J	2233	SER
2	J	2244	LYS
2	J	2247	ASN
2	J	2249	LYS
2	J	2258	ARG
2	J	2259	VAL
2	J	2266	ARG
2	J	2268	LEU
2	J	2303	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	532	ASN
1	B	549	GLN
1	B	786	HIS
1	B	877	GLN
1	B	1517	ASN
1	B	1548	HIS
1	B	1615	ASN
1	B	1667	GLN
1	B	1857	ASN
1	B	1909	GLN
1	B	2016	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](#)

10 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	Q9I	B	2201	-	14,14,14	2.69	7 (50%)	18,20,20	3.42	4 (22%)
4	EDO	B	2208	-	3,3,3	0.47	0	2,2,2	0.39	0
4	EDO	B	2204	-	3,3,3	0.51	0	2,2,2	0.35	0
4	EDO	B	2210	-	3,3,3	0.50	0	2,2,2	0.50	0
4	EDO	B	2203	-	3,3,3	0.55	0	2,2,2	0.57	0
4	EDO	B	2202	-	3,3,3	0.61	0	2,2,2	0.38	0
4	EDO	B	2205	-	3,3,3	0.45	0	2,2,2	1.04	0
4	EDO	B	2207	-	3,3,3	0.46	0	2,2,2	0.42	0
4	EDO	B	2209	-	3,3,3	0.39	0	2,2,2	0.74	0
4	EDO	B	2206	-	3,3,3	0.45	0	2,2,2	0.66	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	Q9I	B	2201	-	-	7/12/14/14	0/1/1/1
4	EDO	B	2208	-	-	1/1/1/1	-
4	EDO	B	2204	-	-	1/1/1/1	-
4	EDO	B	2210	-	-	0/1/1/1	-
4	EDO	B	2203	-	-	1/1/1/1	-
4	EDO	B	2202	-	-	0/1/1/1	-
4	EDO	B	2205	-	-	1/1/1/1	-
4	EDO	B	2207	-	-	1/1/1/1	-
4	EDO	B	2209	-	-	1/1/1/1	-
4	EDO	B	2206	-	-	1/1/1/1	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2201	Q9I	O07-S05	6.13	1.50	1.43
3	B	2201	Q9I	O06-S05	4.62	1.48	1.43
3	B	2201	Q9I	C08-S05	3.13	1.80	1.76
3	B	2201	Q9I	O03-N02	2.67	1.47	1.42
3	B	2201	Q9I	C13-C08	-2.40	1.35	1.38
3	B	2201	Q9I	O03-C04	-2.22	1.39	1.43
3	B	2201	Q9I	C11-CL14	2.01	1.79	1.74

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	2201	Q9I	O07-S05-O06	-12.69	99.78	119.59
3	B	2201	Q9I	O06-S05-C08	4.56	113.77	108.10
3	B	2201	Q9I	O03-N02-C01	2.90	118.90	114.33
3	B	2201	Q9I	C09-C08-S05	-2.56	117.19	119.73

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	2201	Q9I	C13-C08-S05-N02
3	B	2201	Q9I	C09-C08-S05-N02
3	B	2201	Q9I	C01-N02-O03-C04
4	B	2204	EDO	O1-C1-C2-O2
3	B	2201	Q9I	C09-C08-S05-O06
3	B	2201	Q9I	C13-C08-S05-O06
4	B	2209	EDO	O1-C1-C2-O2
3	B	2201	Q9I	C01-N02-S05-C08

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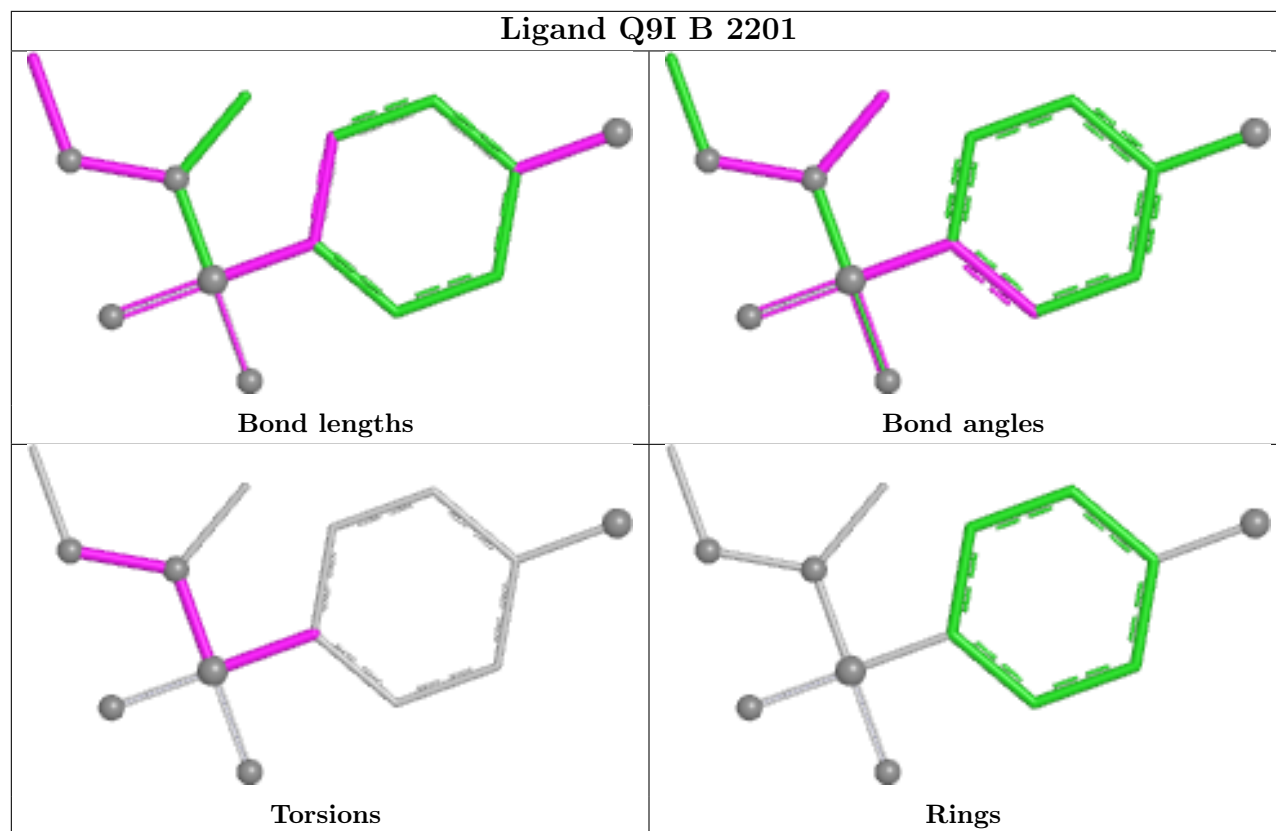
Mol	Chain	Res	Type	Atoms
4	B	2208	EDO	O1-C1-C2-O2
4	B	2206	EDO	O1-C1-C2-O2
4	B	2207	EDO	O1-C1-C2-O2
3	B	2201	Q9I	C01-N02-S05-O07
4	B	2203	EDO	O1-C1-C2-O2
4	B	2205	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2201	Q9I	1	0
4	B	2206	EDO	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	B	1725/1747 (98%)	0.33	82 (4%) 35 32	31, 61, 125, 178	0
2	J	260/263 (98%)	0.14	8 (3%) 51 47	37, 56, 107, 164	0
All	All	1985/2010 (98%)	0.31	90 (4%) 38 34	31, 60, 124, 178	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1895	LEU	5.9
1	B	1859	PRO	4.8
1	B	2038	LEU	4.1
1	B	1849	ILE	4.0
1	B	1889	VAL	3.9
1	B	1940	LEU	3.8
1	B	1836	LEU	3.7
2	J	2320	LEU	3.7
1	B	2035	VAL	3.6
1	B	404	ALA	3.6
1	B	2066	VAL	3.6
1	B	2036	VAL	3.6
1	B	2108	PHE	3.5
1	B	1908	LEU	3.4
1	B	2106	LEU	3.4
1	B	455	PHE	3.3
1	B	456	GLY	3.3
1	B	1911	ASP	3.3
1	B	1868	LEU	3.2
1	B	1858	ILE	3.2
1	B	575	LEU	3.2
2	J	2097	ILE	3.2
1	B	1842	VAL	3.2
1	B	1912	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	2101	ALA	3.1
1	B	1891	THR	3.0
1	B	772	LEU	3.0
1	B	601	GLY	3.0
2	J	2061	GLY	2.9
1	B	1936	LEU	2.9
1	B	698	ILE	2.9
2	J	2063	MET	2.9
1	B	2122	PHE	2.9
1	B	1584	ILE	2.9
1	B	1971	ILE	2.8
1	B	1896	GLN	2.8
1	B	1894	LEU	2.8
1	B	732	GLY	2.8
1	B	1879	LEU	2.8
1	B	1904	LEU	2.8
1	B	1860	ILE	2.7
1	B	778	LEU	2.7
1	B	2107	TYR	2.7
1	B	2037	VAL	2.7
1	B	1044	VAL	2.6
2	J	2062	SER	2.5
1	B	454	PRO	2.4
1	B	1995	ALA	2.4
1	B	1835	SER	2.4
1	B	1982	VAL	2.4
1	B	1597	LEU	2.4
1	B	2068	ILE	2.3
1	B	1852	ALA	2.3
1	B	1419	LEU	2.3
1	B	1601	LEU	2.3
1	B	1893	LEU	2.3
1	B	572	ASP	2.3
2	J	2064	THR	2.3
1	B	1604	LEU	2.3
1	B	1850	SER	2.3
1	B	1845	LEU	2.2
1	B	2024	VAL	2.3
1	B	2039	VAL	2.3
1	B	2104	TYR	2.2
1	B	1871	LEU	2.2
1	B	1075	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	1886	ASP	2.2
1	B	2027	ASP	2.2
1	B	2071	ALA	2.2
1	B	780	TYR	2.2
1	B	1037	LEU	2.2
1	B	1153	LEU	2.2
1	B	573	HIS	2.2
1	B	1978	GLY	2.2
1	B	403	LEU	2.1
1	B	405	PRO	2.1
1	B	859	PRO	2.1
1	B	1855	TYR	2.1
1	B	2090	VAL	2.1
1	B	766	ALA	2.1
1	B	1853	ALA	2.1
1	B	1880	ASN	2.1
2	J	2318	ALA	2.1
1	B	779	PRO	2.0
1	B	1754	TYR	2.0
1	B	782	PHE	2.0
1	B	741	MET	2.0
1	B	1857	ASN	2.0
2	J	2319	LEU	2.0
1	B	2077	ILE	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](#)

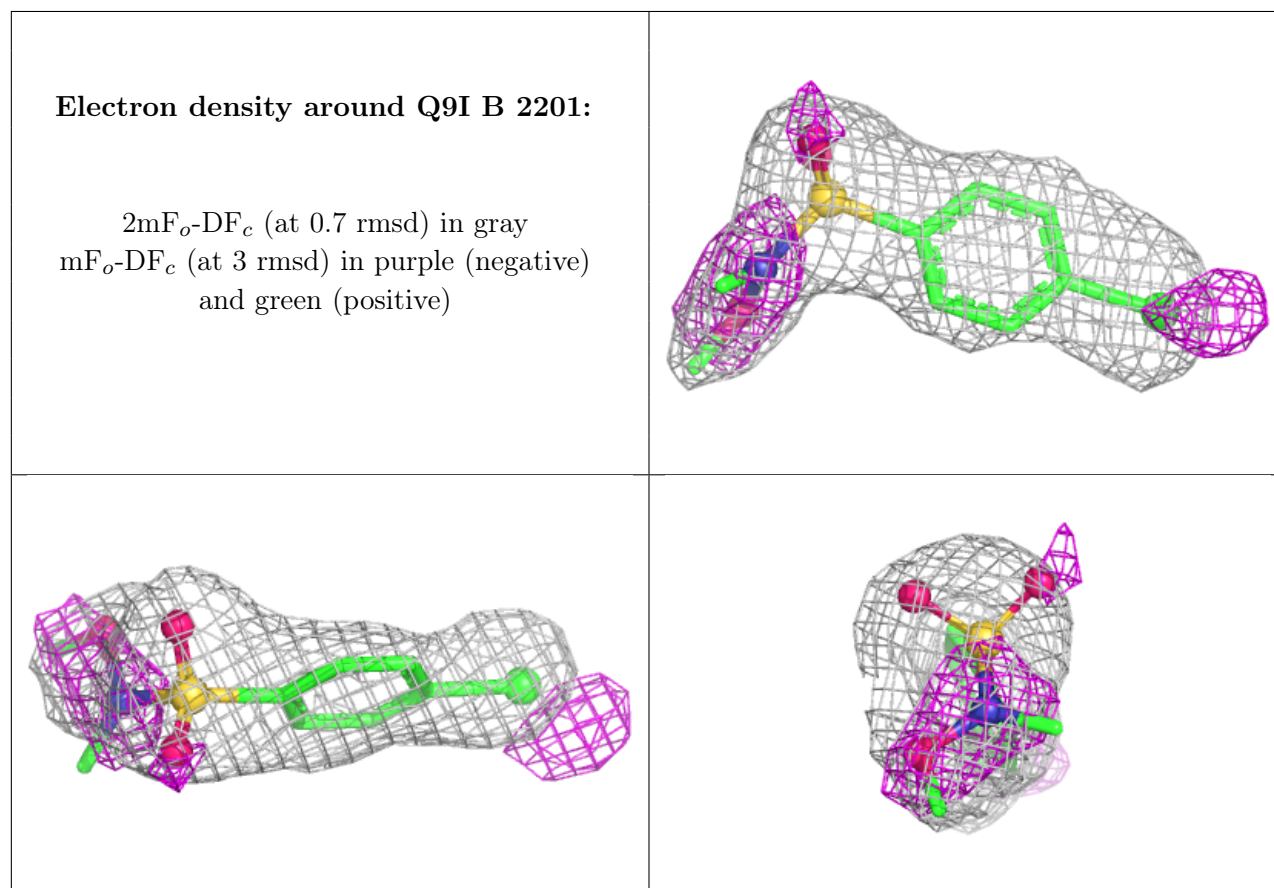
There are no oligosaccharides in this entry.

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	EDO	B	2209	4/4	0.66	0.15	58,59,61,65	0
4	EDO	B	2210	4/4	0.78	0.14	48,51,52,61	0
4	EDO	B	2202	4/4	0.84	0.19	52,54,54,59	0
4	EDO	B	2204	4/4	0.86	0.11	51,54,57,63	0
4	EDO	B	2207	4/4	0.87	0.15	65,67,68,69	0
4	EDO	B	2203	4/4	0.88	0.12	53,54,54,60	0
4	EDO	B	2208	4/4	0.91	0.11	46,50,56,57	0
4	EDO	B	2205	4/4	0.92	0.09	45,48,49,56	0
3	Q9I	B	2201	14/14	0.93	0.10	47,55,59,64	0
4	EDO	B	2206	4/4	0.93	0.09	37,38,43,47	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers ⓘ

There are no such residues in this entry.