



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:34 AM UTC

PDB ID : 4JLU / pdb_00004jlu
Title : Crystal structure of BRCA1 BRCT with doubly phosphorylated Abraxas
Authors : Badgujar, D.; Varma, A.K.
Deposited on : 2013-03-13
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

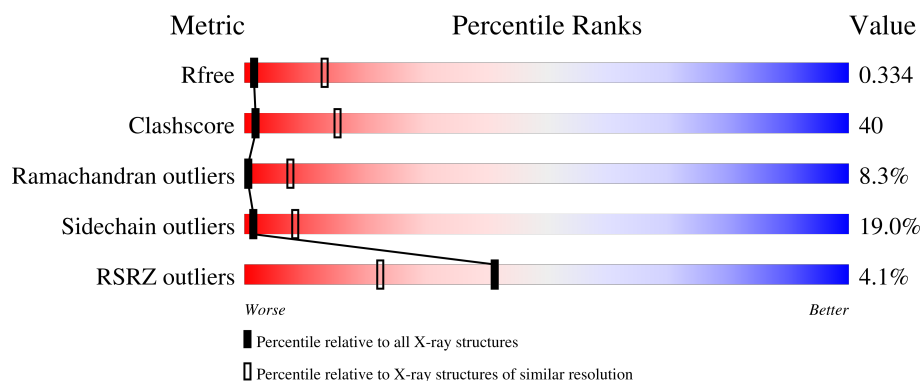
1 Overall quality at a glance

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	211	4% 29% 51% 18% .
2	B	11	9% 18% 55% 18% 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1792 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 Breast cancer type 1 susceptibility protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1696	1082	289	311	14			

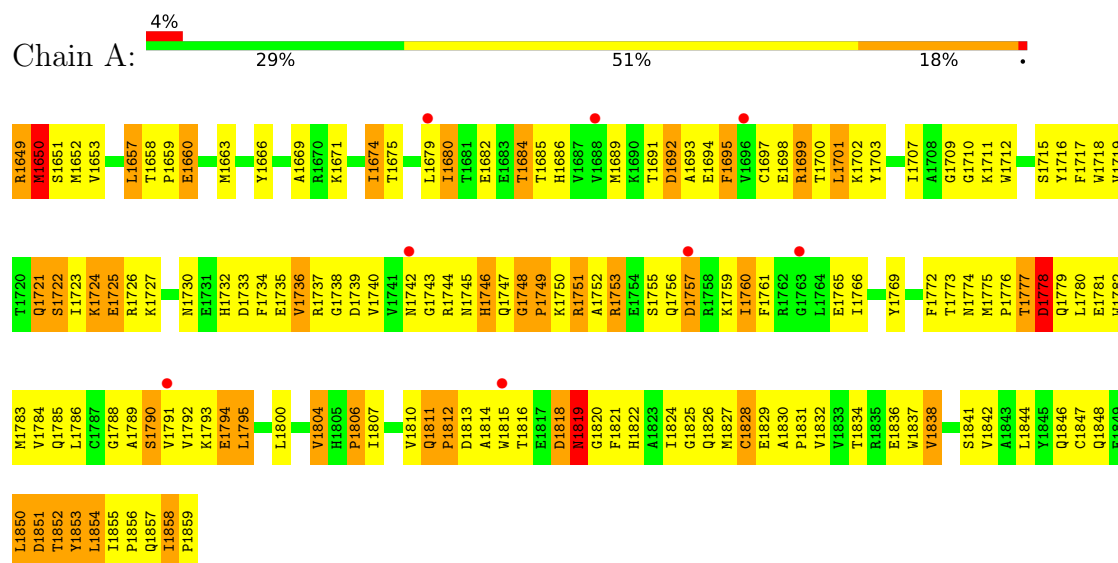
- Molecule 2 is a protein called BRCA1-A complex subunit Abraxas.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	11	Total	C	N	O	P	0	0	0
			96	57	14	23	2			

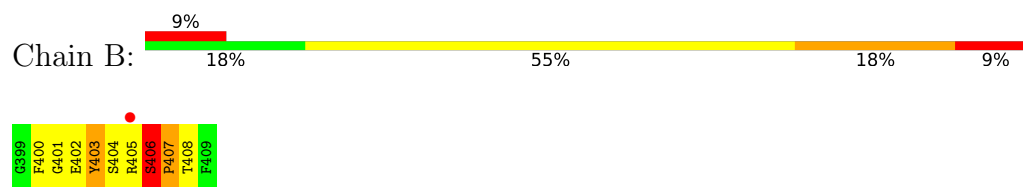
3 Residue-property plots

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: Breast cancer type 1 susceptibility protein



- Molecule 2: BRCA1-A complex subunit Abraxas



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	113.60Å 113.60Å 121.61Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	22.19 – 3.50 22.19 – 3.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (22.19-3.50) 99.4 (22.19-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 3.53Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.298 , 0.327 0.295 , 0.334	Depositor DCC
R_{free} test set	289 reflections (4.65%)	wwPDB-VP
Wilson B-factor (Å ²)	74.3	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 21.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	1792	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.72% 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: SEP

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	0.94	0/1737	1.28	23/2355 (1.0%)
2	B	1.36	0/77	1.25	0/98
All	All	0.96	0/1814	1.28	23/2453 (0.9%)

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
2	B	0	1
All	All	0	2

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1825	GLY	N-CA-C	-10.27	103.38	114.67
1	A	1748	GLY	CA-C-N	8.89	130.95	119.84
1	A	1748	GLY	C-N-CA	8.89	130.95	119.84
1	A	1746	HIS	N-CA-C	8.06	121.09	111.33
1	A	1695	PHE	N-CA-C	7.49	122.14	113.38
1	A	1744	ARG	N-CA-C	-7.02	105.91	114.75
1	A	1650	MET	N-CA-C	6.94	120.04	108.73
1	A	1738	GLY	N-CA-C	6.42	120.19	110.18
1	A	1743	GLY	N-CA-C	6.27	119.37	111.85
1	A	1722	SER	N-CA-C	-6.27	105.67	113.38
1	A	1804	VAL	N-CA-C	5.86	117.21	108.54
1	A	1739	ASP	N-CA-C	5.84	115.81	108.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1680	ILE	N-CA-C	5.80	116.66	108.36
1	A	1818	ASP	N-CA-C	5.68	122.91	110.80
1	A	1830	ALA	CA-C-N	5.67	126.92	119.84
1	A	1830	ALA	C-N-CA	5.67	126.92	119.84
1	A	1742	ASN	N-CA-C	5.57	120.07	112.88
1	A	1760	ILE	N-CA-C	5.55	119.88	111.89
1	A	1666	TYR	N-CA-C	-5.42	105.02	111.03
1	A	1838	VAL	CB-CA-C	-5.34	105.13	111.97
1	A	1657	LEU	N-CA-C	5.14	121.75	110.80
1	A	1658	THR	CA-C-N	5.06	126.16	119.84
1	A	1658	THR	C-N-CA	5.06	126.16	119.84

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1858	ILE	Peptide
2	B	406	SEP	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	1696	0	1671	134	0
2	B	96	0	71	11	0
All	All	1792	0	1742	141	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 40.

All (141) 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:1751:ARG:HG3	1:A:1751:ARG:HH11	1.08	1.11
1:A:1709:GLY:HA3	1:A:1711:LYS:HE2	1.51	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1751:ARG:HG3	1:A:1751:ARG:NH1	1.87	0.88
1:A:1775:MET:HE1	1:A:1783:MET:CE	2.05	0.86
1:A:1700:THR:O	1:A:1703:TYR:HB3	1.76	0.85
2:B:406:SEP:HB3	2:B:407:PRO:HD3	1.60	0.83
1:A:1815:TRP:HB3	1:A:1818:ASP:HB3	1.61	0.81
1:A:1752:ALA:O	1:A:1756:GLN:NE2	2.12	0.81
1:A:1700:THR:HB	2:B:403:TYR:HB2	1.63	0.78
2:B:400:PHE:N	2:B:401:GLY:HA2	2.00	0.76
2:B:400:PHE:H	2:B:401:GLY:HA2	1.49	0.76
2:B:406:SEP:CB	2:B:407:PRO:HD3	2.17	0.74
1:A:1700:THR:HA	2:B:403:TYR:HB3	1.69	0.74
1:A:1751:ARG:HH11	1:A:1751:ARG:CG	1.96	0.74
1:A:1782:TRP:HA	1:A:1785:GLN:HE21	1.53	0.72
1:A:1748:GLY:N	1:A:1749:PRO:HD2	2.04	0.71
1:A:1700:THR:O	1:A:1703:TYR:N	2.20	0.70
1:A:1701:LEU:HD13	1:A:1775:MET:HE3	1.73	0.70
1:A:1674:ILE:HG23	1:A:1675:THR:N	2.07	0.69
1:A:1775:MET:HE1	1:A:1783:MET:HE3	1.75	0.69
1:A:1795:LEU:HD12	1:A:1795:LEU:H	1.57	0.69
1:A:1716:TYR:C	1:A:1718:TRP:H	2.02	0.68
1:A:1782:TRP:HE1	1:A:1786:LEU:HD21	1.59	0.68
1:A:1699:ARG:HA	1:A:1703:TYR:CD2	2.29	0.67
2:B:402:GLU:O	2:B:403:TYR:HD1	1.78	0.67
1:A:1761:PHE:O	1:A:1788:GLY:O	2.12	0.67
1:A:1716:TYR:O	1:A:1718:TRP:N	2.28	0.66
1:A:1834:THR:HG23	1:A:1855:ILE:HD11	1.77	0.66
1:A:1652:MET:CB	1:A:1686:HIS:HB2	2.27	0.65
1:A:1818:ASP:C	1:A:1820:GLY:H	2.05	0.64
1:A:1652:MET:HB3	1:A:1686:HIS:HB2	1.80	0.64
1:A:1815:TRP:HB3	1:A:1818:ASP:CB	2.28	0.63
1:A:1782:TRP:O	1:A:1786:LEU:HG	2.00	0.62
1:A:1819:ASN:HD22	1:A:1822:HIS:HB2	1.65	0.61
1:A:1674:ILE:CG2	1:A:1675:THR:N	2.64	0.60
1:A:1761:PHE:HE2	1:A:1784:VAL:HG12	1.67	0.60
1:A:1781:GLU:HB3	1:A:1791:VAL:HG21	1.84	0.60
1:A:1789:ALA:O	1:A:1790:SER:CB	2.50	0.60
1:A:1848:GLN:HG3	1:A:1853:TYR:OH	2.01	0.59
1:A:1659:PRO:O	1:A:1663:MET:HG2	2.03	0.59
1:A:1700:THR:HA	2:B:403:TYR:CB	2.33	0.58
1:A:1685:THR:O	1:A:1712:TRP:HB2	2.04	0.58
1:A:1748:GLY:H	1:A:1749:PRO:HD2	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1851:ASP:C	1:A:1854:LEU:HD11	2.30	0.56
1:A:1737:ARG:HB3	1:A:1747:GLN:HG3	1.87	0.56
1:A:1811:GLN:HG2	1:A:1812:PRO:HD2	1.85	0.56
1:A:1834:THR:CG2	1:A:1855:ILE:HD11	2.36	0.56
1:A:1836:GLU:O	1:A:1837:TRP:C	2.48	0.56
1:A:1858:ILE:HG22	1:A:1859:PRO:OXT	2.06	0.56
1:A:1775:MET:CE	1:A:1783:MET:HE3	2.34	0.56
1:A:1686:HIS:CD2	1:A:1712:TRP:HB3	2.41	0.56
1:A:1649:ARG:NH2	1:A:1650:MET:HB2	2.21	0.56
1:A:1653:VAL:HG13	1:A:1684:THR:HG21	1.88	0.55
1:A:1699:ARG:HA	1:A:1703:TYR:HD2	1.69	0.55
1:A:1850:LEU:O	1:A:1852:THR:N	2.35	0.54
1:A:1819:ASN:C	1:A:1821:PHE:N	2.65	0.54
1:A:1669:ALA:HB1	1:A:1674:ILE:HG22	1.88	0.54
1:A:1650:MET:HE2	1:A:1674:ILE:HD11	1.90	0.54
1:A:1772:PHE:HD1	1:A:1811:GLN:HB2	1.71	0.54
1:A:1760:ILE:HB	1:A:1847:CYS:HB2	1.89	0.54
1:A:1858:ILE:HG22	1:A:1859:PRO:C	2.33	0.54
1:A:1700:THR:CB	2:B:403:TYR:HB2	2.34	0.53
1:A:1700:THR:O	1:A:1701:LEU:C	2.52	0.53
1:A:1724:LYS:O	1:A:1725:GLU:HG3	2.09	0.53
1:A:1777:THR:O	1:A:1779:GLN:N	2.41	0.53
1:A:1785:GLN:HG3	1:A:1791:VAL:HG23	1.90	0.53
2:B:406:SEP:CB	2:B:407:PRO:CD	2.85	0.53
1:A:1832:VAL:HG21	1:A:1857:GLN:HG3	1.91	0.52
1:A:1674:ILE:HG23	1:A:1675:THR:H	1.73	0.52
1:A:1782:TRP:O	1:A:1782:TRP:CD1	2.63	0.52
1:A:1857:GLN:O	1:A:1859:PRO:HD3	2.10	0.52
1:A:1689:MET:O	1:A:1716:TYR:HB2	2.09	0.52
1:A:1776:PRO:O	1:A:1777:THR:C	2.53	0.51
1:A:1777:THR:O	1:A:1778:ASP:C	2.53	0.51
1:A:1769:TYR:HB3	1:A:1810:VAL:HG12	1.93	0.51
1:A:1700:THR:O	1:A:1703:TYR:CB	2.55	0.51
1:A:1819:ASN:C	1:A:1821:PHE:H	2.18	0.51
1:A:1692:ASP:O	1:A:1694:GLU:N	2.44	0.51
1:A:1781:GLU:OE2	1:A:1791:VAL:HG11	2.11	0.50
1:A:1735:GLU:HG2	1:A:1736:VAL:N	2.25	0.50
1:A:1851:ASP:HA	1:A:1854:LEU:HD21	1.94	0.49
1:A:1850:LEU:C	1:A:1852:THR:H	2.18	0.49
1:A:1652:MET:HB2	1:A:1686:HIS:HB2	1.95	0.49
1:A:1757:ASP:OD1	1:A:1757:ASP:N	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1854:LEU:O	1:A:1855:ILE:C	2.56	0.49
1:A:1701:LEU:CD1	1:A:1775:MET:HE3	2.40	0.49
1:A:1795:LEU:H	1:A:1795:LEU:CD1	2.15	0.48
1:A:1855:ILE:HG23	1:A:1856:PRO:HD2	1.94	0.48
1:A:1719:VAL:C	1:A:1721:GLN:H	2.20	0.48
1:A:1723:ILE:C	1:A:1725:GLU:H	2.21	0.48
1:A:1789:ALA:O	1:A:1790:SER:HB3	2.14	0.48
1:A:1765:GLU:OE1	1:A:1804:VAL:HG11	2.14	0.47
1:A:1679:LEU:HD23	1:A:1702:LYS:HE2	1.98	0.46
1:A:1775:MET:HE1	1:A:1783:MET:HE1	1.91	0.46
1:A:1746:HIS:ND1	1:A:1748:GLY:N	2.64	0.46
1:A:1857:GLN:O	1:A:1859:PRO:CD	2.64	0.46
1:A:1777:THR:O	1:A:1780:LEU:N	2.49	0.45
1:A:1719:VAL:C	1:A:1721:GLN:N	2.75	0.45
1:A:1769:TYR:CE1	1:A:1795:LEU:HD11	2.52	0.45
1:A:1735:GLU:OE2	1:A:1753:ARG:NH2	2.49	0.45
1:A:1822:HIS:HA	1:A:1857:GLN:HE22	1.82	0.45
1:A:1749:PRO:HA	1:A:1752:ALA:HB3	1.99	0.44
1:A:1691:THR:HG21	1:A:1715:SER:OG	2.18	0.44
1:A:1750:LYS:C	1:A:1752:ALA:N	2.73	0.44
1:A:1824:ILE:C	1:A:1826:GLN:N	2.70	0.44
1:A:1811:GLN:NE2	1:A:1813:ASP:OD1	2.50	0.44
1:A:1812:PRO:HD3	1:A:1834:THR:HG22	2.00	0.44
1:A:1701:LEU:C	1:A:1703:TYR:N	2.76	0.44
1:A:1669:ALA:HB1	1:A:1674:ILE:CG2	2.48	0.44
1:A:1651:SER:O	1:A:1652:MET:HB3	2.18	0.44
1:A:1766:ILE:HD12	1:A:1789:ALA:HB2	1.99	0.44
1:A:1806:PRO:O	1:A:1807:ILE:HD13	2.18	0.43
1:A:1773:THR:O	1:A:1774:ASN:HB2	2.18	0.43
1:A:1701:LEU:C	1:A:1703:TYR:H	2.25	0.43
1:A:1777:THR:HG22	1:A:1781:GLU:HG3	2.00	0.43
1:A:1854:LEU:H	1:A:1854:LEU:HD12	1.82	0.43
1:A:1855:ILE:CG2	1:A:1856:PRO:HD2	2.48	0.43
1:A:1707:ILE:C	1:A:1710:GLY:H	2.26	0.43
1:A:1697:CYS:SG	1:A:1698:GLU:N	2.90	0.42
1:A:1748:GLY:C	1:A:1750:LYS:H	2.27	0.42
1:A:1730:ASN:O	1:A:1733:ASP:HB2	2.19	0.42
1:A:1792:VAL:HG12	1:A:1794:GLU:O	2.20	0.42
1:A:1838:VAL:O	1:A:1842:VAL:HG23	2.20	0.42
1:A:1760:ILE:HD12	1:A:1761:PHE:CE1	2.55	0.42
1:A:1682:GLU:H	1:A:1682:GLU:CD	2.28	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1660:GLU:O	1:A:1663:MET:HB2	2.20	0.42
1:A:1811:GLN:NE2	1:A:1814:ALA:HB2	2.34	0.42
1:A:1828:CYS:HB2	1:A:1829:GLU:H	1.70	0.41
1:A:1807:ILE:HD13	1:A:1831:PRO:HG2	2.02	0.41
1:A:1746:HIS:CE1	1:A:1748:GLY:N	2.89	0.41
1:A:1775:MET:HE2	1:A:1780:LEU:HA	2.01	0.41
1:A:1674:ILE:HG23	1:A:1675:THR:O	2.20	0.41
1:A:1837:TRP:O	1:A:1841:SER:OG	2.26	0.41
1:A:1686:HIS:HA	1:A:1712:TRP:O	2.20	0.41
1:A:1766:ILE:HD12	1:A:1789:ALA:CB	2.51	0.41
2:B:400:PHE:N	2:B:401:GLY:CA	2.77	0.41
1:A:1695:PHE:CE2	1:A:1734:PHE:HD1	2.39	0.40
1:A:1775:MET:CE	1:A:1783:MET:CE	2.86	0.40
1:A:1782:TRP:CD1	1:A:1786:LEU:HG	2.56	0.40
1:A:1824:ILE:C	1:A:1827:MET:H	2.29	0.40
1:A:1660:GLU:CD	1:A:1660:GLU:H	2.28	0.40

There are no symmetry-related clashes.

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	209/211 (99%)	140 (67%)	52 (25%)	17 (8%)	1	8
2	B	7/11 (64%)	3 (43%)	3 (43%)	1 (14%)	0	3
All	All	216/222 (97%)	143 (66%)	55 (26%)	18 (8%)	0	7

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1657	LEU
1	A	1693	ALA

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Mol	Chain	Res	Type
1	A	1790	SER
1	A	1819	ASN
1	A	1851	ASP
1	A	1717	PHE
1	A	1725	GLU
1	A	1726	ARG
1	A	1812	PRO
1	A	1816	THR
1	A	1828	CYS
1	A	1749	PRO
1	A	1793	LYS
1	A	1778	ASP
1	A	1671	LYS
1	A	1724	LYS
2	B	407	PRO
1	A	1806	PRO

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	188/188 (100%)	154 (82%)	34 (18%)	2	10
2	B	7/7 (100%)	4 (57%)	3 (43%)	0	0
All	All	195/195 (100%)	158 (81%)	37 (19%)	1	8

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1649	ARG
1	A	1650	MET
1	A	1660	GLU
1	A	1674	ILE
1	A	1680	ILE
1	A	1684	THR
1	A	1692	ASP

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Mol	Chain	Res	Type
1	A	1699	ARG
1	A	1701	LEU
1	A	1721	GLN
1	A	1722	SER
1	A	1727	LYS
1	A	1732	HIS
1	A	1736	VAL
1	A	1740	VAL
1	A	1745	ASN
1	A	1751	ARG
1	A	1753	ARG
1	A	1755	SER
1	A	1757	ASP
1	A	1759	LYS
1	A	1777	THR
1	A	1778	ASP
1	A	1794	GLU
1	A	1795	LEU
1	A	1800	LEU
1	A	1811	GLN
1	A	1819	ASN
1	A	1844	LEU
1	A	1846	GLN
1	A	1850	LEU
1	A	1852	THR
1	A	1853	TYR
1	A	1854	LEU
2	B	403	TYR
2	B	405	ARG
2	B	408	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1672	HIS
1	A	1747	GLN
1	A	1785	GLN
1	A	1819	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SEP	B	404	2	8,9,10	1.89	2 (25%)	7,12,14	2.10	1 (14%)
2	SEP	B	406	2	8,9,10	1.91	3 (37%)	7,12,14	2.05	2 (28%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SEP	B	404	2	-	4/6/8/10	-
2	SEP	B	406	2	-	3/6/8/10	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	404	SEP	P-O1P	3.96	1.62	1.50
2	B	406	SEP	P-O1P	3.81	1.62	1.50
2	B	406	SEP	P-O2P	2.11	1.62	1.54
2	B	406	SEP	P-O3P	2.04	1.62	1.54
2	B	404	SEP	P-O3P	2.04	1.62	1.54

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	404	SEP	OG-CB-CA	4.70	112.72	108.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	406	SEP	OG-CB-CA	4.32	112.35	108.14
2	B	406	SEP	OG-P-O1P	2.59	113.43	106.44

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	404	SEP	N-CA-CB-OG
2	B	406	SEP	CB-OG-P-O1P
2	B	406	SEP	CB-OG-P-O2P
2	B	406	SEP	CB-OG-P-O3P
2	B	404	SEP	CB-OG-P-O2P
2	B	404	SEP	CA-CB-OG-P
2	B	404	SEP	CB-OG-P-O3P

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	406	SEP	3	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	211/211 (100%)	0.55	8 (3%) 44 24	43, 50, 50, 50	0
2	B	9/11 (81%)	1.08	1 (11%) 10 7	25, 50, 50, 50	1 (11%)
All	All	220/222 (99%)	0.57	9 (4%) 41 22	25, 50, 50, 50	1 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	405	ARG	5.1
1	A	1757	ASP	2.8
1	A	1763	GLY	2.6
1	A	1815	TRP	2.4
1	A	1679	LEU	2.3
1	A	1742	ASN	2.3
1	A	1696	VAL	2.3
1	A	1791	VAL	2.2
1	A	1688	VAL	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	SEP	B	404	10/11	0.85	0.13	50,50,50,50	0
2	SEP	B	406	10/11	0.87	0.12	50,50,50,50	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.