



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:00 PM UTC

PDB ID : 2ING / pdb\_00002ing  
Title : X-ray Structure of the BRCA1 BRCT mutant M1775K  
Authors : Birrane, G.; Soni, A.; Ladas, J.A.A.  
Deposited on : 2006-10-07  
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](mailto: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>.

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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

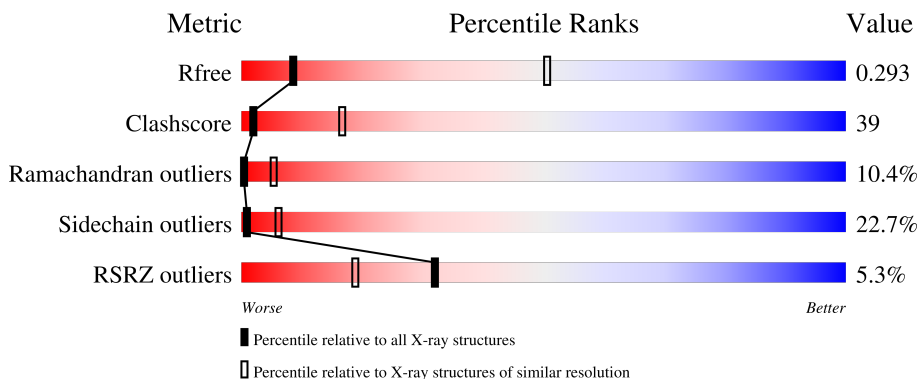
# 1 Overall quality at a glance

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  $\geq 3$ , 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	X	213	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1682 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
			Total	C	N	O	S			
1	X	208	1671	1069	287	302	13	0	0	0

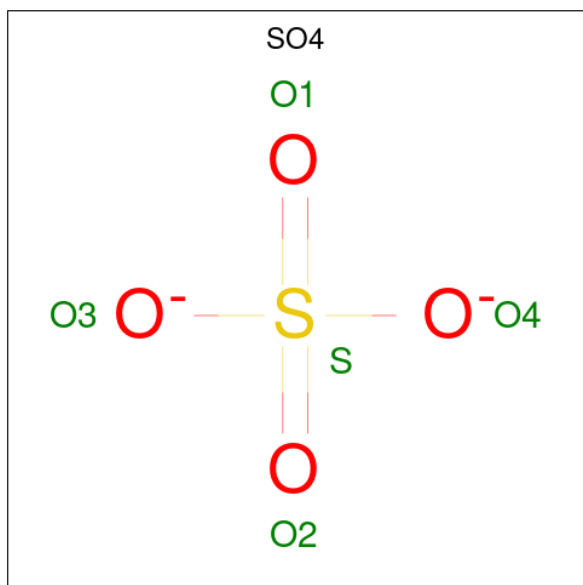
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	1647	GLY	-	expression tag	UNP P38398
X	1648	PRO	-	expression tag	UNP P38398
X	1775	LYS	MET	engineered mutation	UNP P38398

- Molecule 2 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	X	1	Total	Co	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O<sub>4</sub>S).

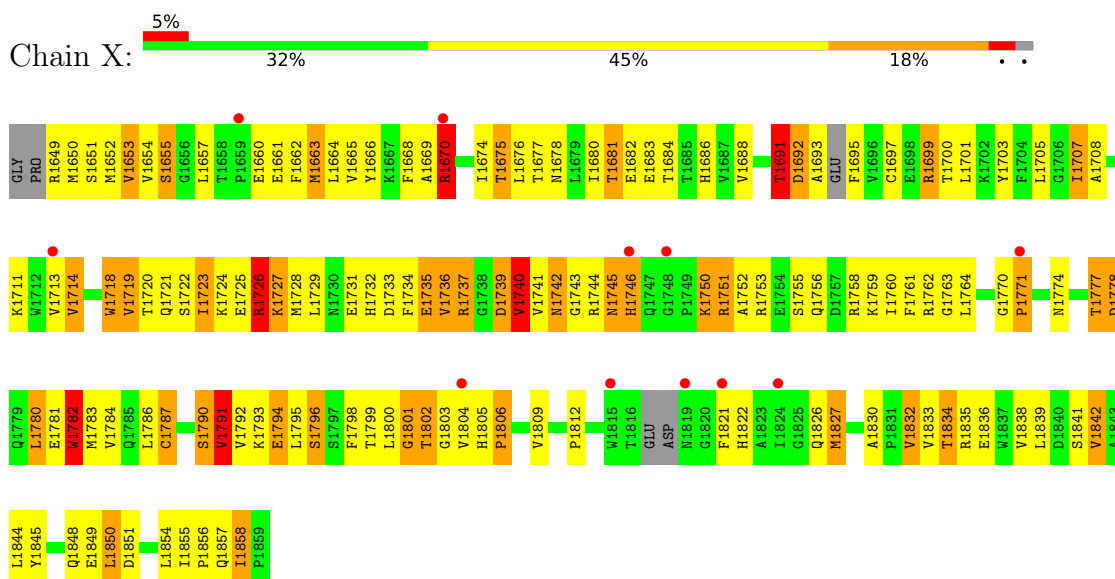


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	X	1	Total	O	S	0	0
			5	4	1		
3	X	1	Total	O	S	0	0
			5	4	1		

### 3 Residue-property plots [i](#)

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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	114.34Å 114.34Å 119.88Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 3.60 30.00 – 3.60	Depositor EDS
% Data completeness (in resolution range)	99.0 (30.00-3.60) 98.7 (30.00-3.60)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	9.96 (at 3.65Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.247 , 0.302 0.244 , 0.293	Depositor DCC
$R_{free}$ test set	501 reflections (8.74%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	136.9	Xtrriage
Anisotropy	0.149	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 254.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.36$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	1682	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	174.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.31% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, CO

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	X	0.86	1/1710 (0.1%)	1.09	7/2315 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	1719	VAL	CA-CB	-7.77	1.46	1.55

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1770	GLY	CA-C-N	7.12	128.74	119.84
1	X	1770	GLY	C-N-CA	7.12	128.74	119.84
1	X	1655	SER	N-CA-C	6.31	118.73	108.26
1	X	1746	HIS	N-CA-C	5.64	118.08	109.62
1	X	1707	ILE	CB-CA-C	-5.20	105.21	112.02
1	X	1791	VAL	N-CA-C	5.18	115.23	107.51
1	X	1782	TRP	CA-CB-CG	5.01	123.11	113.60

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	X	1671	0	1657	129	0
2	X	1	0	0	0	0
3	X	10	0	0	2	0
All	All	1682	0	1657	129	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 39.

All (129) 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:X:1653:VAL:HG13	1:X:1677:THR:O	1.50	1.09
1:X:1653:VAL:HG22	1:X:1677:THR:HG22	1.42	1.00
1:X:1653:VAL:HG22	1:X:1677:THR:CG2	2.01	0.90
1:X:1838:VAL:O	1:X:1842:VAL:HG23	1.74	0.86
1:X:1832:VAL:HG21	1:X:1857:GLN:NE2	1.91	0.85
1:X:1741:VAL:HG12	1:X:1844:LEU:HD11	1.64	0.80
1:X:1835:ARG:HD3	1:X:1839:LEU:HD11	1.63	0.80
1:X:1684:THR:HB	1:X:1711:LYS:HZ3	1.48	0.78
1:X:1652:MET:HE2	1:X:1674:ILE:HG21	1.66	0.76
1:X:1697:CYS:SG	1:X:1736:VAL:HG21	2.26	0.74
1:X:1722:SER:OG	1:X:1729:LEU:HD13	1.91	0.70
1:X:1725:GLU:O	1:X:1727:LYS:N	2.25	0.69
1:X:1801:GLY:O	1:X:1803:GLY:N	2.25	0.69
1:X:1761:PHE:HB2	1:X:1787:CYS:O	1.92	0.69
1:X:1760:ILE:HG23	1:X:1845:TYR:CZ	2.29	0.68
1:X:1794:GLU:CD	1:X:1794:GLU:N	2.54	0.65
1:X:1700:THR:O	1:X:1703:TYR:HB3	1.97	0.65
1:X:1668:PHE:CG	1:X:1719:VAL:HG13	2.31	0.65
1:X:1855:ILE:HG22	1:X:1857:GLN:NE2	2.12	0.65
1:X:1718:TRP:HA	1:X:1734:PHE:CE2	2.33	0.64
1:X:1669:ALA:O	1:X:1670:ARG:C	2.41	0.64
1:X:1681:THR:O	1:X:1711:LYS:NZ	2.32	0.63
1:X:1723:ILE:HG22	1:X:1724:LYS:N	2.14	0.63
1:X:1800:LEU:O	1:X:1801:GLY:C	2.42	0.61
1:X:1692:ASP:O	1:X:1693:ALA:HB3	2.00	0.61
1:X:1653:VAL:CG2	1:X:1677:THR:CG2	2.78	0.59
1:X:1832:VAL:HG21	1:X:1857:GLN:CD	2.27	0.59
1:X:1654:VAL:HG21	1:X:1657:LEU:CD1	2.31	0.59
1:X:1809:VAL:HG22	1:X:1833:VAL:HG23	1.85	0.59
1:X:1841:SER:HA	1:X:1848:GLN:HE21	1.68	0.59
1:X:1668:PHE:CD1	1:X:1719:VAL:HG13	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:1718:TRP:CZ2	1:X:1728:MET:HE2	2.38	0.59
1:X:1742:ASN:HB3	1:X:1746:HIS:HB2	1.85	0.59
1:X:1791:VAL:HG12	1:X:1791:VAL:O	2.02	0.58
1:X:1826:GLN:C	1:X:1827:MET:HE3	2.28	0.58
1:X:1751:ARG:O	1:X:1752:ALA:C	2.46	0.58
1:X:1663:MET:HE3	1:X:1664:LEU:CD2	2.33	0.58
1:X:1741:VAL:CG1	1:X:1844:LEU:HD11	2.33	0.57
1:X:1680:ILE:O	1:X:1680:ILE:HG23	2.04	0.56
1:X:1735:GLU:OE1	1:X:1753:ARG:CZ	2.53	0.56
1:X:1723:ILE:CG2	1:X:1724:LYS:N	2.69	0.56
1:X:1851:ASP:O	1:X:1854:LEU:HD12	2.06	0.55
1:X:1661:GLU:O	1:X:1665:VAL:HG23	2.06	0.55
1:X:1760:ILE:HG21	1:X:1845:TYR:O	2.07	0.55
1:X:1683:GLU:N	1:X:1683:GLU:OE1	2.40	0.54
1:X:1707:ILE:HG23	1:X:1752:ALA:HB2	1.89	0.54
1:X:1822:HIS:O	1:X:1858:ILE:HG22	2.08	0.54
1:X:1651:SER:HA	1:X:1675:THR:HG22	1.89	0.53
1:X:1701:LEU:O	1:X:1705:LEU:HD12	2.09	0.53
1:X:1654:VAL:HG23	1:X:1688:VAL:HB	1.90	0.53
1:X:1777:THR:CG2	1:X:1778:ASP:N	2.71	0.53
1:X:1778:ASP:N	1:X:1778:ASP:OD1	2.39	0.53
1:X:1782:TRP:O	1:X:1786:LEU:HD13	2.09	0.53
1:X:1854:LEU:C	1:X:1855:ILE:HD12	2.33	0.53
1:X:1695:PHE:CB	1:X:1737:ARG:HD3	2.39	0.53
1:X:1834:THR:O	1:X:1835:ARG:C	2.51	0.53
1:X:1678:ASN:ND2	3:X:2:SO4:O4	2.42	0.52
1:X:1792:VAL:HG21	1:X:1798:PHE:CD2	2.45	0.52
1:X:1761:PHE:CE1	1:X:1784:VAL:HG12	2.45	0.52
1:X:1653:VAL:HG12	1:X:1654:VAL:H	1.76	0.51
1:X:1697:CYS:O	1:X:1739:ASP:HB3	2.11	0.51
1:X:1705:LEU:HA	1:X:1708:ALA:HB3	1.91	0.50
1:X:1800:LEU:HD13	1:X:1802:THR:N	2.27	0.50
1:X:1664:LEU:HD12	1:X:1720:THR:HG22	1.92	0.50
1:X:1654:VAL:HG21	1:X:1657:LEU:HD11	1.92	0.50
1:X:1850:LEU:HD23	1:X:1850:LEU:N	2.26	0.50
1:X:1692:ASP:O	1:X:1693:ALA:CB	2.60	0.50
1:X:1855:ILE:HG23	1:X:1856:PRO:HD2	1.93	0.50
1:X:1806:PRO:HG2	1:X:1830:ALA:HB2	1.94	0.49
1:X:1699:ARG:NE	1:X:1741:VAL:CG2	2.75	0.49
1:X:1707:ILE:O	1:X:1708:ALA:C	2.55	0.49
1:X:1664:LEU:O	1:X:1668:PHE:N	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:1654:VAL:HG21	1:X:1657:LEU:HD12	1.94	0.49
1:X:1782:TRP:CD1	1:X:1786:LEU:HD13	2.48	0.49
1:X:1781:GLU:CD	1:X:1793:LYS:HZ1	2.21	0.49
1:X:1682:GLU:OE2	1:X:1782:TRP:CH2	2.65	0.48
1:X:1799:THR:O	1:X:1800:LEU:C	2.56	0.48
1:X:1663:MET:HE3	1:X:1664:LEU:HD21	1.93	0.48
1:X:1652:MET:CE	1:X:1674:ILE:HG21	2.41	0.48
1:X:1654:VAL:CG2	1:X:1657:LEU:CD1	2.92	0.48
1:X:1740:VAL:HG12	1:X:1741:VAL:N	2.28	0.48
1:X:1660:GLU:O	1:X:1663:MET:HB3	2.14	0.47
1:X:1699:ARG:NE	1:X:1741:VAL:HG23	2.28	0.47
1:X:1756:GLN:HE22	1:X:1759:LYS:HZ1	1.62	0.47
1:X:1750:LYS:O	1:X:1751:ARG:C	2.57	0.47
1:X:1688:VAL:HA	1:X:1714:VAL:O	2.14	0.46
1:X:1700:THR:HG21	3:X:1:SO4:O3	2.15	0.46
1:X:1725:GLU:O	1:X:1726:ARG:C	2.58	0.46
1:X:1855:ILE:HG22	1:X:1857:GLN:HE22	1.81	0.46
1:X:1652:MET:HB3	1:X:1676:LEU:HD12	1.97	0.46
1:X:1674:ILE:HG22	1:X:1675:THR:O	2.16	0.46
1:X:1844:LEU:H	1:X:1844:LEU:HD12	1.80	0.46
1:X:1700:THR:HG22	1:X:1701:LEU:N	2.31	0.46
1:X:1763:GLY:O	1:X:1764:LEU:HD23	2.16	0.46
1:X:1718:TRP:HA	1:X:1734:PHE:CZ	2.51	0.45
1:X:1662:PHE:O	1:X:1663:MET:C	2.60	0.45
1:X:1699:ARG:HA	1:X:1703:TYR:CD2	2.52	0.45
1:X:1654:VAL:CG2	1:X:1657:LEU:HD11	2.47	0.44
1:X:1794:GLU:O	1:X:1795:LEU:C	2.59	0.44
1:X:1763:GLY:C	1:X:1764:LEU:HD23	2.43	0.44
1:X:1783:MET:O	1:X:1786:LEU:HB2	2.18	0.44
1:X:1822:HIS:C	1:X:1858:ILE:HG22	2.43	0.44
1:X:1680:ILE:O	1:X:1680:ILE:CG2	2.66	0.43
1:X:1691:THR:HG22	1:X:1695:PHE:C	2.42	0.43
1:X:1695:PHE:HB2	1:X:1737:ARG:HH21	1.82	0.43
1:X:1850:LEU:O	1:X:1851:ASP:C	2.60	0.43
1:X:1663:MET:HE3	1:X:1664:LEU:HD23	1.99	0.43
1:X:1699:ARG:HA	1:X:1703:TYR:HD2	1.84	0.43
1:X:1731:GLU:C	1:X:1733:ASP:H	2.27	0.43
1:X:1780:LEU:O	1:X:1783:MET:HB3	2.18	0.43
1:X:1662:PHE:C	1:X:1662:PHE:CD2	2.97	0.43
1:X:1663:MET:O	1:X:1666:TYR:HB3	2.19	0.43
1:X:1841:SER:HA	1:X:1848:GLN:NE2	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:1650:MET:HE3	1:X:1686:HIS:CE1	2.53	0.43
1:X:1760:ILE:HG23	1:X:1845:TYR:CE1	2.54	0.43
1:X:1858:ILE:O	1:X:1858:ILE:HG23	2.19	0.42
1:X:1700:THR:O	1:X:1701:LEU:C	2.61	0.42
1:X:1662:PHE:O	1:X:1662:PHE:HD2	2.02	0.42
1:X:1735:GLU:OE1	1:X:1753:ARG:NE	2.53	0.42
1:X:1674:ILE:HG22	1:X:1675:THR:N	2.34	0.42
1:X:1742:ASN:OD1	1:X:1742:ASN:N	2.51	0.41
1:X:1850:LEU:O	1:X:1854:LEU:CD1	2.68	0.41
1:X:1662:PHE:C	1:X:1662:PHE:HD2	2.29	0.41
1:X:1783:MET:HE3	1:X:1783:MET:HB2	1.85	0.41
1:X:1764:LEU:O	1:X:1790:SER:HB2	2.21	0.41
1:X:1850:LEU:C	1:X:1854:LEU:HD11	2.46	0.41
1:X:1742:ASN:ND2	1:X:1746:HIS:O	2.50	0.41
1:X:1805:HIS:HA	1:X:1806:PRO:HD3	1.87	0.41
1:X:1737:ARG:O	1:X:1743:GLY:O	2.40	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	X	202/213 (95%)	148 (73%)	33 (16%)	21 (10%)	<b>0</b> <b>5</b>

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	1726	ARG
1	X	1740	VAL
1	X	1750	LYS
1	X	1771	PRO
1	X	1802	THR

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Mol	Chain	Res	Type
1	X	1692	ASP
1	X	1739	ASP
1	X	1745	ASN
1	X	1762	ARG
1	X	1663	MET
1	X	1821	PHE
1	X	1732	HIS
1	X	1755	SER
1	X	1670	ARG
1	X	1691	THR
1	X	1718	TRP
1	X	1806	PRO
1	X	1836	GLU
1	X	1796	SER
1	X	1801	GLY
1	X	1812	PRO

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	X	185/189 (98%)	143 (77%)	42 (23%)	<b>1</b> <b>6</b>

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	1649	ARG
1	X	1653	VAL
1	X	1655	SER
1	X	1670	ARG
1	X	1675	THR
1	X	1681	THR
1	X	1691	THR
1	X	1699	ARG
1	X	1713	VAL
1	X	1714	VAL

*Continued on next page...*

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	X	1721	GLN
1	X	1723	ILE
1	X	1726	ARG
1	X	1727	LYS
1	X	1735	GLU
1	X	1736	VAL
1	X	1737	ARG
1	X	1740	VAL
1	X	1742	ASN
1	X	1744	ARG
1	X	1745	ASN
1	X	1751	ARG
1	X	1758	ARG
1	X	1771	PRO
1	X	1774	ASN
1	X	1777	THR
1	X	1778	ASP
1	X	1780	LEU
1	X	1782	TRP
1	X	1787	CYS
1	X	1790	SER
1	X	1791	VAL
1	X	1794	GLU
1	X	1796	SER
1	X	1804	VAL
1	X	1827	MET
1	X	1832	VAL
1	X	1834	THR
1	X	1842	VAL
1	X	1849	GLU
1	X	1850	LEU
1	X	1858	ILE

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

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	X	1756	GLN
1	X	1774	ASN
1	X	1779	GLN
1	X	1819	ASN
1	X	1848	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](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	SO4	X	1	-	4,4,4	0.37	0	6,6,6	0.36	0
3	SO4	X	2	-	4,4,4	0.22	0	6,6,6	0.41	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	X	1	SO4	1	0
3	X	2	SO4	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	X	208/213 (97%)	0.76	11 (5%) 32 19	175, 175, 175, 175	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	X	1815	TRP	4.6
1	X	1771	PRO	3.1
1	X	1746	HIS	3.0
1	X	1748	GLY	2.3
1	X	1804	VAL	2.3
1	X	1819	ASN	2.3
1	X	1824	ILE	2.2
1	X	1670	ARG	2.1
1	X	1659	PRO	2.1
1	X	1821	PHE	2.1
1	X	1713	VAL	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	SO4	X	1	5/5	0.77	0.32	174,174,174,174	0
3	SO4	X	2	5/5	0.80	0.29	174,174,174,174	0
2	CO	X	101	1/1	0.89	0.11	174,174,174,174	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.