



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:15 AM UTC

PDB ID : 4ZZZ / pdb_00004zzz
Title : Structure of human PARP1 catalytic domain bound to an isoindolinone inhibitor
Authors : Casale, E.; Fasolini, M.; Papeo, G.; Posterl, H.; Borghi, D.; Busel, A.A.; Caprera, F.; Ciomei, M.; Cirila, A.; Corti, E.; DAnello, M.; Fasolini, M.; Felder, E.R.; Forte, B.; Galvani, A.; Isacchi, A.; Khvat, A.; Krasavin, M.Y.; Lupi, R.; Orsini, P.; Perego, R.; Pesenti, E.; Pezzetta, D.; Rainoldi, S.; RiccardiSirtori, F.; Scolaro, A.; Sola, F.; Zuccotto, F.; Donati, D.; Montagnoli, A.
Deposited on : 2015-04-15
Resolution : 1.90 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)

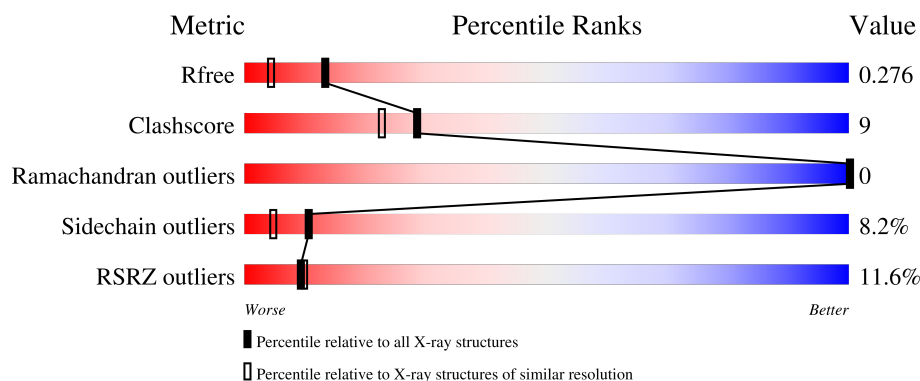
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	362	
1	B	362	

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5676 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 POLY [ADP-RIBOSE] POLYMERASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	350	Total	C	N	O	S	0	0	0
			2761	1762	466	522	11			
1	B	342	Total	C	N	O	S	0	0	0
			2695	1720	455	509	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	653	GLY	-	expression tag	UNP P09874
A	654	PRO	-	expression tag	UNP P09874
B	653	GLY	-	expression tag	UNP P09874
B	654	PRO	-	expression tag	UNP P09874

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



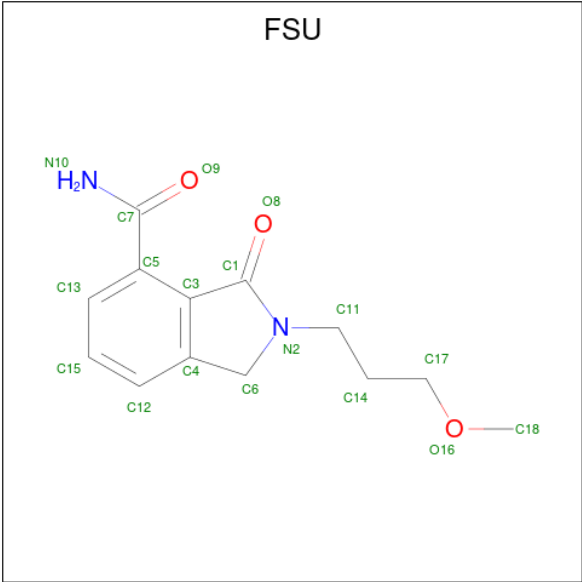
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

- Molecule 4 is 2-(3-methoxypropyl)-3-oxo-2,3-dihydro-1H-isoindole-4-carboxamide (CCD ID: FSU) (formula: $C_{13}H_{16}N_2O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			18	13	2	3		
4	B	1	Total	C	N	O	0	0
			18	13	2	3		

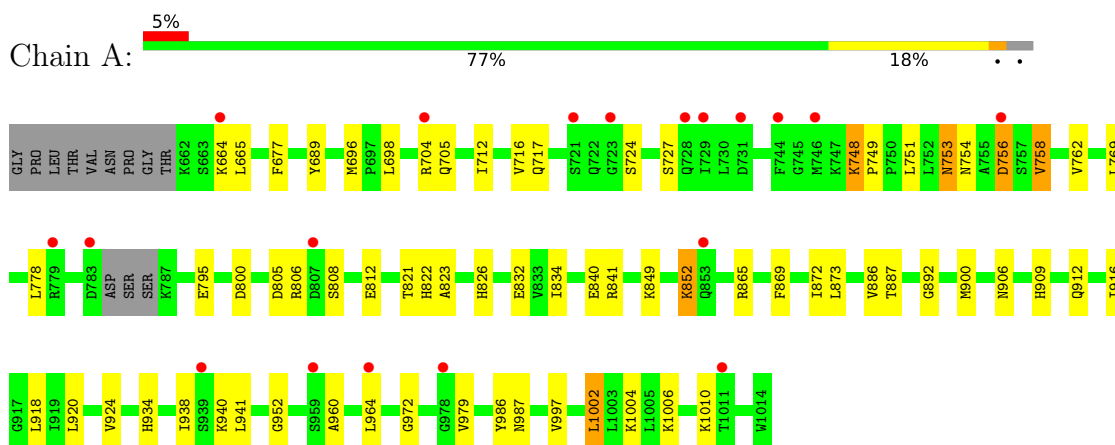
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	79	Total	O	0	0
			79	79		
5	B	67	Total	O	0	0
			67	67		

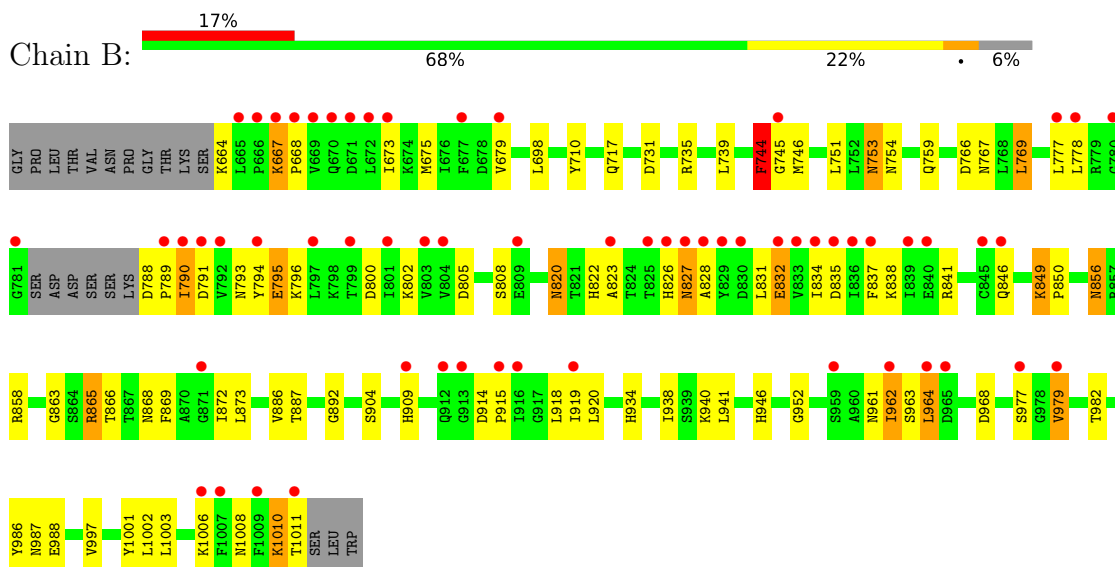
3 Residue-property plots

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: POLY [ADP-RIBOSE] POLYMERASE 1



• Molecule 1: POLY [ADP-RIBOSE] POLYMERASE 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	179.90Å 54.05Å 92.76Å 90.00° 113.79° 90.00°	Depositor
Resolution (Å)	30.00 – 1.90 30.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	96.7 (30.00-1.90) 96.7 (30.00-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.228 , 0.265 (Not available) , 0.276	Depositor DCC
R_{free} test set	3161 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.459	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 37.6	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	0.94	EDS
Total number of atoms	5676	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.76% 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: SO4, FSU, GOL

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	0.85	0/2814	0.96	1/3798 (0.0%)
1	B	0.80	1/2746 (0.0%)	0.98	4/3707 (0.1%)
All	All	0.83	1/5560 (0.0%)	0.97	5/7505 (0.1%)

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	946	HIS	ND1-CE1	5.05	1.37	1.32

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	744	PHE	N-CA-C	6.29	120.95	113.16
1	B	869	PHE	N-CA-C	6.11	118.90	111.82
1	A	979	VAL	N-CA-C	5.83	116.50	107.99
1	B	1010	LYS	N-CA-C	-5.58	102.64	110.50
1	B	739	LEU	N-CA-C	5.36	116.81	111.07

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	744	PHE	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	2761	0	2806	38	0
1	B	2695	0	2740	64	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	12	0	16	0	0
3	B	6	0	8	0	0
4	A	18	0	16	0	0
4	B	18	0	16	1	0
5	A	79	0	0	2	0
5	B	67	0	0	8	0
All	All	5676	0	5602	98	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 9.

All (98) 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:1003:LEU:HD12	5:B:3067:HOH:O	1.45	1.11
1:B:793:ASN:HA	1:B:796:LYS:HE2	1.51	0.91
1:B:769:LEU:HD22	5:B:3016:HOH:O	1.83	0.77
1:B:796:LYS:HD3	5:B:3018:HOH:O	1.86	0.74
1:A:841:ARG:HD2	1:A:873:LEU:O	1.88	0.73
1:B:667:LYS:CG	1:B:668:PRO:HD3	2.19	0.73
1:B:962:ILE:HG22	1:B:964:LEU:HD22	1.70	0.72
1:A:952:GLY:H	1:A:987:ASN:HD22	1.38	0.72
1:B:1010:LYS:O	1:B:1011:THR:HB	1.92	0.70
1:B:667:LYS:HG3	1:B:668:PRO:HD3	1.75	0.69
1:B:919:ILE:HG22	1:B:1003:LEU:HB2	1.75	0.67
1:B:1001:TYR:HB3	5:B:3067:HOH:O	1.94	0.66
1:B:769:LEU:CD2	5:B:3016:HOH:O	2.41	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:822:HIS:HE1	1:A:826:HIS:O	1.82	0.62
1:B:952:GLY:H	1:B:987:ASN:ND2	1.98	0.62
1:A:952:GLY:H	1:A:987:ASN:ND2	1.98	0.61
1:A:886:VAL:HG11	1:B:934:HIS:CD2	2.35	0.61
1:A:717:GLN:HG2	1:A:887:THR:OG1	2.00	0.60
1:B:952:GLY:H	1:B:987:ASN:HD22	1.49	0.58
1:B:961:ASN:HD21	1:B:968:ASP:HB3	1.68	0.58
1:B:1006:LYS:HE2	1:B:1008:ASN:ND2	2.19	0.58
1:B:872:ILE:HG21	1:B:920:LEU:HD11	1.87	0.57
1:A:886:VAL:CG1	1:B:934:HIS:CD2	2.88	0.57
1:B:962:ILE:CG2	1:B:964:LEU:HD22	2.35	0.56
1:B:778:LEU:HA	1:B:796:LYS:HZ2	1.72	0.55
1:A:805:ASP:O	1:A:808:SER:OG	2.23	0.55
1:B:822:HIS:HE1	1:B:826:HIS:O	1.91	0.54
1:B:834:ILE:HD11	1:B:1006:LYS:HB2	1.90	0.54
1:A:689:TYR:OH	1:A:865:ARG:NH1	2.42	0.53
1:A:872:ILE:HG21	1:A:920:LEU:HD11	1.89	0.53
1:B:805:ASP:HB2	1:B:808:SER:HB3	1.89	0.53
1:A:834:ILE:HG21	1:A:1004:LYS:HE2	1.90	0.52
1:A:916:ILE:HD11	1:A:1006:LYS:HD2	1.91	0.52
1:B:777:LEU:HD22	1:B:796:LYS:HB3	1.89	0.52
1:A:918:LEU:HD22	1:A:1002:LEU:HD21	1.92	0.52
1:B:710:TYR:OH	1:B:766:ASP:OD1	2.24	0.52
1:B:915:PRO:HD2	5:B:3052:HOH:O	2.10	0.52
1:B:827:ASN:HD22	1:B:828:ALA:N	2.07	0.52
1:B:865:ARG:NH1	1:B:909:HIS:ND1	2.58	0.52
1:A:849:LYS:O	1:A:852:LYS:HB2	2.10	0.52
1:B:832:GLU:HG2	1:B:1006:LYS:HB3	1.91	0.51
1:B:918:LEU:HD22	1:B:1002:LEU:HD21	1.91	0.51
1:B:841:ARG:HD2	1:B:873:LEU:O	2.09	0.51
1:B:753:ASN:ND2	1:B:754:ASN:HD22	2.08	0.51
1:B:667:LYS:HG2	1:B:668:PRO:HD3	1.92	0.50
1:B:793:ASN:HA	1:B:796:LYS:CE	2.31	0.50
1:B:820:ASN:HD22	1:B:820:ASN:N	2.11	0.48
1:A:696:MET:SD	1:A:705:GLN:NE2	2.76	0.48
1:B:794:TYR:O	1:B:795:GLU:C	2.56	0.48
1:A:758:VAL:O	1:A:762:VAL:HG23	2.14	0.47
1:B:835:ASP:HB2	1:B:837:PHE:CE1	2.49	0.47
1:A:677:PHE:HB3	1:A:778:LEU:HD13	1.97	0.47
1:A:712:ILE:O	1:A:716:VAL:HG23	2.15	0.46
1:A:892:GLY:HA3	1:A:938:ILE:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:796:LYS:CD	5:B:3018:HOH:O	2.56	0.46
1:A:960:ALA:HB3	1:A:972:GLY:CA	2.45	0.46
1:A:822:HIS:HD2	5:A:3026:HOH:O	1.98	0.46
1:B:753:ASN:HD22	1:B:753:ASN:H	1.64	0.45
1:A:869:PHE:HE2	1:A:920:LEU:HG	1.81	0.45
1:B:731:ASP:OD2	1:B:735:ARG:NH1	2.50	0.45
1:B:1010:LYS:O	1:B:1011:THR:CB	2.60	0.45
1:A:823:ALA:HB1	1:A:986:TYR:CZ	2.50	0.45
1:A:832:GLU:HG3	1:A:1006:LYS:HB3	1.99	0.45
1:A:916:ILE:CD1	1:A:1006:LYS:HD2	2.48	0.44
1:B:767:ASN:HB3	5:B:3015:HOH:O	2.17	0.44
1:A:821:THR:HB	1:A:900:MET:HA	2.00	0.44
1:B:667:LYS:HG3	1:B:668:PRO:CD	2.47	0.44
1:A:753:ASN:HD22	1:A:753:ASN:C	2.26	0.43
1:B:673:ILE:HD11	1:B:794:TYR:HD1	1.82	0.43
1:B:863:GLY:HA3	1:B:904:SER:O	2.18	0.43
1:A:934:HIS:CD2	1:B:886:VAL:CG1	3.01	0.43
1:B:892:GLY:HA3	1:B:938:ILE:O	2.18	0.43
1:A:1010:LYS:HD2	5:A:3025:HOH:O	2.18	0.43
1:B:790:ILE:HG22	1:B:791:ASP:OD1	2.18	0.43
1:B:849:LYS:HB3	1:B:850:PRO:CD	2.49	0.43
1:A:754:ASN:O	1:A:758:VAL:HG13	2.19	0.42
1:B:977:SER:C	1:B:979:VAL:H	2.28	0.42
1:A:826:HIS:CD2	1:A:906:ASN:HD21	2.38	0.42
1:A:865:ARG:NH1	1:A:909:HIS:ND1	2.67	0.42
1:B:675:MET:HG2	1:B:866:THR:HG21	2.02	0.41
1:B:831:LEU:HA	1:B:1006:LYS:O	2.20	0.41
1:B:919:ILE:CG2	1:B:1003:LEU:HB2	2.48	0.41
1:A:748:LYS:HA	1:A:749:PRO:HD3	1.97	0.41
1:B:823:ALA:HB1	1:B:986:TYR:CZ	2.56	0.41
1:B:856:ASN:OD1	1:B:858:ARG:NH2	2.53	0.41
1:B:865:ARG:O	1:B:868:ASN:HB2	2.21	0.41
1:A:934:HIS:CD2	1:B:886:VAL:HG11	2.55	0.41
1:B:788:ASP:HA	1:B:789:PRO:HD2	1.86	0.40
1:B:717:GLN:HG2	1:B:887:THR:OG1	2.20	0.40
1:A:800:ASP:HB3	1:A:840:GLU:HG3	2.03	0.40
1:A:960:ALA:HB3	1:A:972:GLY:HA3	2.04	0.40
1:B:753:ASN:HD22	1:B:753:ASN:N	2.19	0.40
1:B:988:GLU:HG3	4:B:2015:FSU:H12	2.04	0.40
1:A:753:ASN:HD22	1:A:754:ASN:HD22	1.69	0.40
1:A:754:ASN:OD1	1:A:756:ASP:HB2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:679:VAL:HG13	1:B:698:LEU:HD23	2.03	0.40
1:B:914:ASP:OD1	1:B:914:ASP:C	2.62	0.40
1:B:744:PHE:N	1:B:745:GLY:HA2	2.36	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	A	50/362 (14%)	49 (98%)	1 (2%)	0	100	100
1	B	19/362 (5%)	19 (100%)	0	0	100	100
All	All	69/724 (10%)	68 (99%)	1 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	A	308/318 (97%)	285 (92%)	23 (8%)	12	6
1	B	300/318 (94%)	273 (91%)	27 (9%)	9	3
All	All	608/636 (96%)	558 (92%)	50 (8%)	10	5

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	664	LYS
1	A	665	LEU
1	A	698	LEU
1	A	704	ARG
1	A	724	SER
1	A	727	SER
1	A	748	LYS
1	A	751	LEU
1	A	753	ASN
1	A	756	ASP
1	A	758	VAL
1	A	769	LEU
1	A	795	GLU
1	A	806	ARG
1	A	812	GLU
1	A	852	LYS
1	A	912	GLN
1	A	924	VAL
1	A	940	LYS
1	A	941	LEU
1	A	964	LEU
1	A	997	VAL
1	A	1002	LEU
1	B	664	LYS
1	B	667	LYS
1	B	746	MET
1	B	751	LEU
1	B	753	ASN
1	B	759	GLN
1	B	769	LEU
1	B	790	ILE
1	B	795	GLU
1	B	800	ASP
1	B	802	LYS
1	B	820	ASN
1	B	827	ASN
1	B	832	GLU
1	B	838	LYS
1	B	846	GLN
1	B	849	LYS
1	B	856	ASN
1	B	865	ARG
1	B	940	LYS

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Mol	Chain	Res	Type
1	B	941	LEU
1	B	962	ILE
1	B	963	SER
1	B	964	LEU
1	B	979	VAL
1	B	982	THR
1	B	997	VAL

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

Mol	Chain	Res	Type
1	A	717	GLN
1	A	753	ASN
1	A	820	ASN
1	A	822	HIS
1	A	875	GLN
1	A	906	ASN
1	A	987	ASN
1	A	998	ASN
1	B	670	GLN
1	B	717	GLN
1	B	722	GLN
1	B	753	ASN
1	B	767	ASN
1	B	820	ASN
1	B	822	HIS
1	B	826	HIS
1	B	827	ASN
1	B	868	ASN
1	B	961	ASN
1	B	987	ASN
1	B	1008	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

9 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	FSU	B	2015	-	19,19,19	4.38	8 (42%)	23,26,26	3.59	10 (43%)
3	GOL	A	2018	-	5,5,5	0.18	0	5,5,5	1.16	1 (20%)
2	SO4	B	2014	-	4,4,4	0.42	0	6,6,6	0.20	0
3	GOL	A	2017	-	5,5,5	0.26	0	5,5,5	0.72	0
4	FSU	A	2019	-	19,19,19	4.22	7 (36%)	23,26,26	3.34	7 (30%)
3	GOL	B	2013	-	5,5,5	0.45	0	5,5,5	0.24	0
2	SO4	A	2015	-	4,4,4	0.42	0	6,6,6	0.37	0
2	SO4	A	2016	-	4,4,4	0.47	0	6,6,6	0.23	0
2	SO4	B	2012	-	4,4,4	0.50	0	6,6,6	0.58	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	FSU	B	2015	-	-	2/9/21/21	0/2/2/2
3	GOL	A	2018	-	-	1/4/4/4	-
3	GOL	A	2017	-	-	2/4/4/4	-
4	FSU	A	2019	-	-	3/9/21/21	0/2/2/2
3	GOL	B	2013	-	-	4/4/4/4	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	2015	FSU	C6-C4	-11.08	1.36	1.50
4	A	2019	FSU	C6-C4	-10.98	1.36	1.50
4	B	2015	FSU	C6-N2	-8.98	1.34	1.46
4	A	2019	FSU	C6-N2	-8.27	1.35	1.46
4	B	2015	FSU	C3-C1	-8.23	1.34	1.49
4	A	2019	FSU	C3-C1	-7.69	1.35	1.49
4	A	2019	FSU	O8-C1	6.14	1.34	1.22
4	A	2019	FSU	C5-C7	-5.45	1.42	1.50
4	B	2015	FSU	O8-C1	5.41	1.33	1.22
4	B	2015	FSU	C5-C7	-4.38	1.44	1.50
4	B	2015	FSU	C12-C4	-3.99	1.33	1.39
4	B	2015	FSU	C5-C3	-3.16	1.36	1.41
4	B	2015	FSU	C7-N10	2.42	1.37	1.33
4	A	2019	FSU	C11-N2	2.30	1.53	1.47
4	A	2019	FSU	C12-C4	-2.06	1.36	1.39

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2019	FSU	C4-C6-N2	11.79	107.96	102.42
4	B	2015	FSU	C4-C6-N2	11.54	107.84	102.42
4	A	2019	FSU	C3-C1-N2	7.67	111.95	106.40
4	B	2015	FSU	C3-C1-N2	6.82	111.34	106.40
4	B	2015	FSU	C6-N2-C11	5.94	129.62	123.19
4	B	2015	FSU	O9-C7-C5	-4.03	115.48	120.24
4	A	2019	FSU	C6-N2-C1	-3.68	106.42	113.14
4	B	2015	FSU	C3-C5-C7	3.68	129.24	122.17
4	B	2015	FSU	C5-C7-N10	3.31	123.43	118.29
4	B	2015	FSU	C6-N2-C1	-3.27	107.17	113.14
4	A	2019	FSU	C3-C5-C7	3.26	128.44	122.17
4	A	2019	FSU	O8-C1-N2	-3.01	119.06	125.26
4	B	2015	FSU	C13-C15-C12	-2.83	116.60	120.24
4	A	2019	FSU	C6-N2-C11	2.53	125.93	123.19
4	B	2015	FSU	C13-C5-C3	-2.38	116.80	120.16
3	A	2018	GOL	C3-C2-C1	-2.33	103.26	111.80
4	B	2015	FSU	C13-C5-C7	-2.08	113.88	118.78
4	A	2019	FSU	C13-C5-C7	-2.00	114.06	118.78

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2017	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	B	2013	GOL	O1-C1-C2-C3
4	B	2015	FSU	C3-C5-C7-O9
4	B	2015	FSU	C3-C5-C7-N10
3	B	2013	GOL	C1-C2-C3-O3
3	A	2017	GOL	O1-C1-C2-O2
3	B	2013	GOL	O1-C1-C2-O2
4	A	2019	FSU	C14-C17-O16-C18
4	A	2019	FSU	C3-C5-C7-O9
4	A	2019	FSU	C3-C5-C7-N10
3	B	2013	GOL	O2-C2-C3-O3
3	A	2018	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2015	FSU	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 ⓘ

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	350/362 (96%)	0.47	19 (5%) 31 33	12, 28, 49, 57	0
1	B	342/362 (94%)	0.95	61 (17%) 4 3	15, 34, 52, 72	0
All	All	692/724 (95%)	0.71	80 (11%) 9 10	12, 31, 51, 72	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	781	GLY	5.7
1	B	665	LEU	4.5
1	B	1011	THR	4.4
1	B	789	PRO	4.2
1	B	669	VAL	4.0
1	B	979	VAL	3.7
1	B	834	ILE	3.5
1	B	668	PRO	3.5
1	A	978	GLY	3.5
1	A	723	GLY	3.4
1	A	746	MET	3.4
1	B	673	ILE	3.3
1	A	964	LEU	3.2
1	B	836	ILE	3.2
1	B	677	PHE	3.2
1	B	837	PHE	3.2
1	B	845	CYS	3.2
1	B	809	GLU	3.1
1	B	790	ILE	3.1
1	B	792	VAL	3.1
1	B	871	GLY	3.1
1	B	823	ALA	3.1
1	B	828	ALA	3.0
1	A	1011	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	791	ASP	3.0
1	B	835	ASP	2.9
1	B	666	PRO	2.8
1	A	807	ASP	2.8
1	B	919	ILE	2.8
1	A	729	ILE	2.8
1	A	783	ASP	2.7
1	B	916	ILE	2.7
1	B	1006	LYS	2.7
1	B	804	VAL	2.7
1	A	728	GLN	2.6
1	A	744	PHE	2.6
1	A	721	SER	2.6
1	B	833	VAL	2.6
1	B	801	ILE	2.6
1	B	827	ASN	2.6
1	B	909	HIS	2.5
1	B	778	LEU	2.5
1	B	964	LEU	2.5
1	B	780	GLY	2.5
1	B	777	LEU	2.5
1	B	913	GLY	2.5
1	A	853	GLN	2.5
1	B	745	GLY	2.4
1	A	731	ASP	2.4
1	B	803	VAL	2.4
1	A	939	SER	2.4
1	B	794	TYR	2.4
1	B	1009	PHE	2.4
1	B	840	GLU	2.4
1	B	846	GLN	2.4
1	B	1007	PHE	2.4
1	B	962	ILE	2.4
1	B	826	HIS	2.3
1	B	915	PRO	2.3
1	B	965	ASP	2.3
1	B	797	LEU	2.3
1	B	829	TYR	2.3
1	B	839	ILE	2.2
1	A	664	LYS	2.2
1	B	672	LEU	2.2
1	B	912	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	959	SER	2.2
1	B	667	LYS	2.2
1	B	670	GLN	2.2
1	B	799	THR	2.2
1	B	825	THR	2.2
1	A	756	ASP	2.2
1	B	830	ASP	2.1
1	A	959	SER	2.1
1	B	977	SER	2.1
1	B	832	GLU	2.1
1	B	679	VAL	2.1
1	A	779	ARG	2.1
1	B	671	ASP	2.0
1	A	704	ARG	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
2	SO4	B	2014	5/5	0.83	0.10	46,51,52,55	0
3	GOL	A	2017	6/6	0.92	0.11	23,35,38,43	0
3	GOL	B	2013	6/6	0.92	0.10	26,32,34,37	0
4	FSU	B	2015	18/18	0.92	0.10	18,20,33,33	0
3	GOL	A	2018	6/6	0.93	0.09	23,29,30,37	0
2	SO4	A	2016	5/5	0.94	0.11	43,45,48,49	0
4	FSU	A	2019	18/18	0.96	0.06	13,15,30,31	0
2	SO4	A	2015	5/5	0.98	0.06	19,21,23,25	0
2	SO4	B	2012	5/5	0.98	0.06	26,29,29,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.