



Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2024 – 09:32 PM EDT

PDB ID : 6MFZ
Title : Crystal structure of dimodular LgrA in a condensation state
Authors : Reimer, J.M.; Eivaskhani, M.; Harb, I.; Schmeing, T.M.
Deposited on : 2018-09-12
Resolution : 6.00 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	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

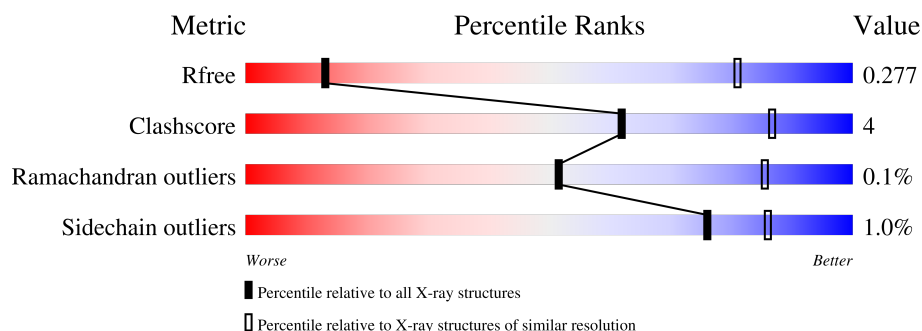
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 6.00 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}	130704	1000 (8.00-3.88)
Clashscore	141614	1049 (8.00-3.90)
Ramachandran outliers	138981	1016 (8.00-3.86)
Sidechain outliers	138945	1017 (8.00-3.82)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	1814	 88% 11% .
1	B	1814	 29% 68% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 18882 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 Linear gramicidin synthase subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1789	Total	C	N	O	S	0	0	0
			14226	9089	2430	2649	58			
1	B	585	Total	C	N	O	S	0	0	0
			4648	2958	797	871	22			

There are 26 discrepancies between the modelled and reference sequences:

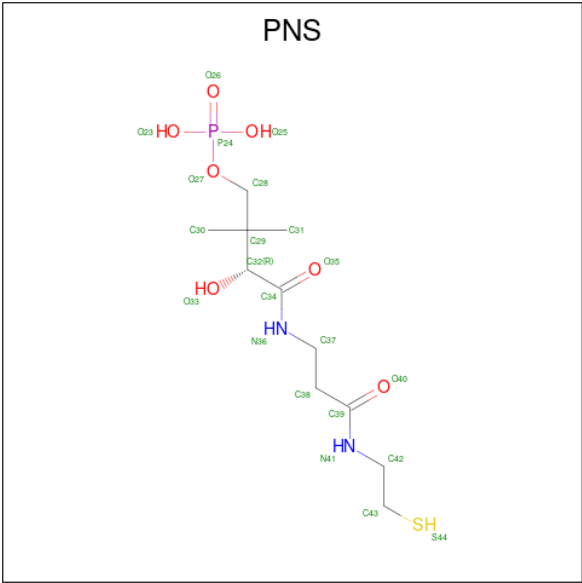
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q70LM7
A	0	ALA	-	expression tag	UNP Q70LM7
A	1	MET	-	expression tag	UNP Q70LM7
A	2	GLY	-	expression tag	UNP Q70LM7
A	1804	ALA	-	expression tag	UNP Q70LM7
A	1805	ALA	-	expression tag	UNP Q70LM7
A	1806	ALA	-	expression tag	UNP Q70LM7
A	1807	GLU	-	expression tag	UNP Q70LM7
A	1808	ASN	-	expression tag	UNP Q70LM7
A	1809	LEU	-	expression tag	UNP Q70LM7
A	1810	TYR	-	expression tag	UNP Q70LM7
A	1811	PHE	-	expression tag	UNP Q70LM7
A	1812	GLN	-	expression tag	UNP Q70LM7
B	-1	GLY	-	expression tag	UNP Q70LM7
B	0	ALA	-	expression tag	UNP Q70LM7
B	1	MET	-	expression tag	UNP Q70LM7
B	2	GLY	-	expression tag	UNP Q70LM7
B	1804	ALA	-	expression tag	UNP Q70LM7
B	1805	ALA	-	expression tag	UNP Q70LM7
B	1806	ALA	-	expression tag	UNP Q70LM7
B	1807	GLU	-	expression tag	UNP Q70LM7
B	1808	ASN	-	expression tag	UNP Q70LM7
B	1809	LEU	-	expression tag	UNP Q70LM7
B	1810	TYR	-	expression tag	UNP Q70LM7
B	1811	PHE	-	expression tag	UNP Q70LM7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1812	GLN	-	expression tag	UNP Q70LM7

- Molecule 2 is 4'-PHOSPHOPANTETHEINE (three-letter code: PNS) (formula: C₁₁H₂₃N₂O₇PS).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			4	3	1		
2	A	1	Total	O	P	0	0
			4	3	1		

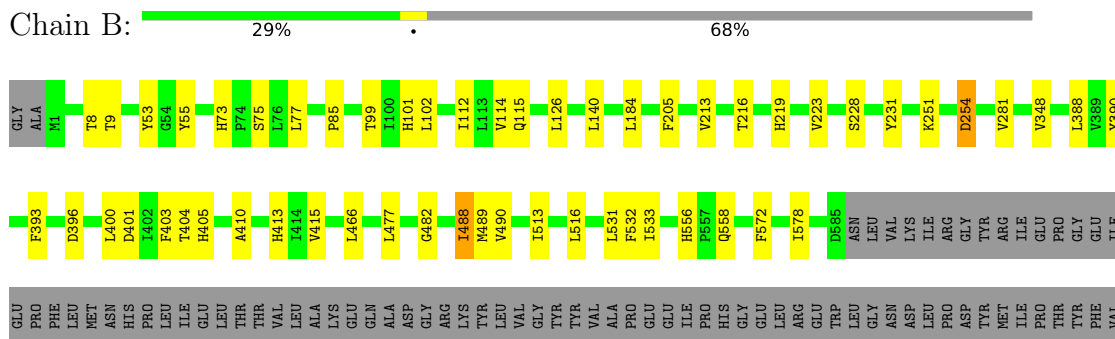
3 Residue-property plots

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. 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. 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: Linear gramicidin synthase subunit A



- Molecule 1: Linear gramicidin synthase subunit A





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	213.60Å 262.75Å 249.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	78.63 – 6.00 78.63 – 6.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (78.63-6.00) 99.8 (78.63-6.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 6.18Å)	Xtriage
Refinement program	PHENIX (dev_3494: ???)	Depositor
R, R_{free}	0.255 , 0.279 0.256 , 0.277	Depositor DCC
R_{free} test set	894 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	216.9	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 455.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	18882	wwPDB-VP
Average B, all atoms (Å ²)	343.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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: PNS

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	0.29	0/14537	0.43	0/19739
1	B	0.25	0/4754	0.40	0/6451
All	All	0.28	0/19291	0.42	0/26190

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	A	14226	0	14159	136	0
1	B	4648	0	4573	30	0
2	A	8	0	0	0	0
All	All	18882	0	18732	166	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 4.

All (166) 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:1531:SER:HB2	1:A:1532:PRO:HD3	1.35	1.07
1:A:1601:LEU:O	1:A:1613:LEU:HB3	1.55	1.05
1:A:407:LEU:HD23	1:A:407:LEU:O	1.61	1.00
1:A:437:ILE:HG23	1:A:458:VAL:HG12	1.42	0.99
1:A:437:ILE:HG12	1:A:458:VAL:HB	1.44	0.96
1:A:368:LEU:HD21	1:A:403:PHE:CD2	2.07	0.89
1:A:1223:ILE:HG23	1:A:1424:VAL:HG11	1.53	0.88
1:A:1295:GLY:HA2	1:A:1358:ASN:OD1	1.75	0.85
1:A:1752:ASN:HB3	1:A:1755:GLU:HB2	1.58	0.85
1:A:1642:ILE:HD13	1:A:1663:LEU:HD22	1.58	0.84
1:A:24:LEU:O	1:A:26:HIS:HD2	1.61	0.83
1:A:1424:VAL:HB	1:A:1425:PRO:HD3	1.64	0.79
1:A:990:ILE:HG12	1:A:1160:GLU:HG2	1.64	0.79
1:A:448:PHE:CZ	1:A:456:PHE:HE2	2.03	0.77
1:A:437:ILE:HG23	1:A:458:VAL:CG1	2.15	0.77
1:A:800:ASP:HB3	1:A:803:SER:HB2	1.68	0.76
1:A:24:LEU:O	1:A:26:HIS:CD2	2.39	0.75
1:A:1664:LEU:HD13	1:A:1696:GLN:HA	1.70	0.74
1:A:1026:LEU:O	1:A:1030:ASN:ND2	2.20	0.70
1:A:995:VAL:HB	1:A:1155:LEU:HB2	1.74	0.68
1:B:126:LEU:HD21	1:B:184:LEU:HD22	1.76	0.68
1:A:1424:VAL:HB	1:A:1425:PRO:CD	2.25	0.67
1:A:402:ILE:O	1:A:406:LEU:HB2	1.95	0.67
1:A:368:LEU:HD21	1:A:403:PHE:CE2	2.31	0.66
1:A:1785:TYR:HB3	1:A:1790:GLN:CB	2.26	0.66
1:A:695:PRO:HB3	1:A:718:ILE:HD11	1.80	0.64
1:A:126:LEU:HD21	1:A:184:LEU:HD22	1.80	0.63
1:A:1111:HIS:HB3	1:A:1112:PRO:HD2	1.81	0.63
1:A:448:PHE:CZ	1:A:456:PHE:CE2	2.86	0.63
1:A:1531:SER:CB	1:A:1532:PRO:HD3	2.18	0.62
1:A:396:ASP:HB3	1:A:488:ILE:HG12	1.80	0.62
1:A:1368:TYR:HA	1:A:1378:GLY:HA2	1.81	0.62
1:A:990:ILE:CG1	1:A:1160:GLU:HG2	2.30	0.60
1:B:513:ILE:CD1	1:B:533:ILE:HG12	2.31	0.60
1:A:1785:TYR:HB3	1:A:1790:GLN:HB3	1.83	0.60
1:A:201:ILE:HG22	1:A:203:LYS:H	1.67	0.60
1:A:552:LYS:O	1:A:554:ILE:HG23	2.02	0.60
1:A:795:ILE:HG12	1:A:1764:LEU:HD22	1.84	0.60
1:B:400:LEU:O	1:B:404:THR:OG1	2.18	0.59
1:A:1337:VAL:O	1:A:1337:VAL:HG12	2.03	0.59
1:A:437:ILE:CG1	1:A:458:VAL:HB	2.25	0.58
1:B:396:ASP:HB2	1:B:488:ILE:HG12	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1773:ILE:O	1:A:1773:ILE:HG22	2.05	0.57
1:B:513:ILE:HD12	1:B:531:LEU:HD21	1.87	0.57
1:A:1069:HIS:HA	1:A:1193:VAL:HB	1.86	0.57
1:A:1464:ILE:HB	1:A:1465:PRO:HD3	1.87	0.57
1:A:1645:ALA:HA	1:A:1663:LEU:HA	1.86	0.56
1:A:1644:ASP:O	1:A:1664:LEU:N	2.36	0.56
1:A:1223:ILE:CG2	1:A:1424:VAL:HG11	2.32	0.56
1:A:1508:ASN:HD22	1:A:1533:VAL:HG12	1.70	0.56
1:A:223:VAL:HG13	1:A:228:SER:HB3	1.87	0.56
1:A:1531:SER:HB2	1:A:1532:PRO:CD	2.23	0.56
1:A:407:LEU:O	1:A:407:LEU:CD2	2.47	0.56
1:A:1664:LEU:HD11	1:A:1697:PHE:H	1.70	0.56
1:A:436:THR:OG1	1:A:437:ILE:HD12	2.07	0.55
1:B:223:VAL:HG13	1:B:228:SER:HB3	1.89	0.55
1:A:458:VAL:HG12	1:A:458:VAL:O	2.06	0.54
1:A:732:VAL:O	1:A:735:LEU:HB3	2.07	0.54
1:A:1329:VAL:HG21	1:A:1340:ASP:OD1	2.07	0.54
1:A:633:ALA:O	1:A:661:HIS:NE2	2.38	0.54
1:A:8:THR:OG1	1:A:9:THR:N	2.41	0.54
1:B:75:SER:HB3	1:B:85:PRO:HB2	1.90	0.54
1:A:1775:TRP:CZ3	1:A:1795:VAL:O	2.62	0.53
1:A:393:PHE:HA	1:A:398:SER:OG	2.08	0.53
1:A:482:GLY:HA3	1:A:489:MET:HA	1.89	0.53
1:A:1664:LEU:CD1	1:A:1696:GLN:HA	2.37	0.53
1:A:1419:GLU:OE2	1:A:1457:ASN:HB2	2.09	0.53
1:A:1295:GLY:HA2	1:A:1358:ASN:CG	2.28	0.52
1:A:281:VAL:HG22	1:A:348:VAL:HB	1.91	0.52
1:A:1145:THR:HB	1:A:1160:GLU:HB2	1.91	0.52
1:A:75:SER:HB3	1:A:85:PRO:HB2	1.90	0.52
1:A:1423:TRP:CD1	1:A:1430:LEU:HB2	2.44	0.52
1:A:668:THR:OG1	1:A:670:ASN:OD1	2.28	0.51
1:B:281:VAL:HG22	1:B:348:VAL:HB	1.93	0.51
1:A:1248:THR:OG1	1:A:1249:TYR:N	2.44	0.51
1:A:1280:ARG:HG2	1:A:1432:ILE:HG23	1.92	0.51
1:A:600:ILE:HD12	1:A:615:VAL:HG11	1.93	0.50
1:B:251:LYS:O	1:B:254:ASP:HB2	2.11	0.50
1:B:482:GLY:HA3	1:B:489:MET:HA	1.94	0.50
1:B:101:HIS:ND1	1:B:102:LEU:O	2.42	0.50
1:B:488:ILE:HG22	1:B:489:MET:HG3	1.94	0.50
1:A:448:PHE:CE2	1:A:456:PHE:CE2	3.00	0.49
1:A:1359:GLN:HG3	1:A:1361:GLY:H	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:990:ILE:HG23	1:A:1158:VAL:HG13	1.94	0.49
1:A:1177:GLN:HE22	1:A:1555:PRO:HB3	1.77	0.49
1:B:8:THR:OG1	1:B:9:THR:N	2.46	0.49
1:B:466:LEU:HD21	1:B:477:LEU:HD21	1.94	0.49
1:A:488:ILE:HG22	1:A:489:MET:HG3	1.95	0.49
1:A:1457:ASN:OD1	1:A:1483:LEU:HD12	2.13	0.49
1:B:405:HIS:HB3	1:B:410:ALA:HB3	1.94	0.49
1:A:600:ILE:HB	1:A:615:VAL:HG21	1.94	0.48
1:A:1158:VAL:HG12	1:A:1160:GLU:HG3	1.95	0.48
1:A:400:LEU:O	1:A:404:THR:CB	2.61	0.48
1:A:388:LEU:HD23	1:A:413:HIS:HB2	1.95	0.48
1:A:125:THR:HA	1:A:180:THR:HA	1.95	0.48
1:A:1295:GLY:HA2	1:A:1358:ASN:ND2	2.28	0.48
1:A:1764:LEU:HG	1:A:1764:LEU:O	2.14	0.48
1:A:1785:TYR:HB3	1:A:1790:GLN:HB2	1.95	0.48
1:A:1538:PRO:HB3	1:A:1544:MET:HG3	1.94	0.48
1:A:1642:ILE:CD1	1:A:1663:LEU:HD22	2.38	0.48
1:B:114:VAL:HG11	1:B:140:LEU:HD11	1.96	0.48
1:A:1414:ASP:O	1:A:1517:ILE:HG12	2.13	0.47
1:A:805:ALA:HA	1:A:1139:ILE:HG21	1.96	0.47
1:A:1560:GLY:O	1:A:1601:LEU:HD12	2.14	0.47
1:A:490:VAL:HG11	1:A:533:ILE:HD13	1.97	0.47
1:B:388:LEU:HD23	1:B:413:HIS:HB2	1.96	0.47
1:B:390:TYR:HA	1:B:415:VAL:HG11	1.96	0.47
1:B:556:HIS:CE1	1:B:558:GLN:HB2	2.50	0.47
1:A:1663:LEU:O	1:A:1663:LEU:HD12	2.14	0.47
1:A:990:ILE:CD1	1:A:1160:GLU:HG2	2.45	0.46
1:A:1277:MET:HB3	1:A:1301:ILE:HB	1.98	0.46
1:A:1561:GLU:HA	1:A:1600:ASP:O	2.15	0.46
1:A:516:LEU:HB2	1:A:532:PHE:CE1	2.51	0.46
1:A:989:ALA:HB3	1:A:1161:TYR:CZ	2.51	0.46
1:A:1301:ILE:O	1:A:1411:TYR:OH	2.22	0.46
1:A:1737:ILE:HD13	1:A:1770:LEU:HD23	1.97	0.45
1:A:390:TYR:HA	1:A:415:VAL:HG11	1.98	0.45
1:A:453:ASN:CG	1:A:456:PHE:CD2	2.90	0.45
1:A:448:PHE:HZ	1:A:456:PHE:HE2	1.60	0.45
1:A:794:TYR:HB3	1:A:848:ILE:HD13	1.97	0.45
1:A:1295:GLY:HA2	1:A:1358:ASN:HD21	1.81	0.45
1:A:101:HIS:ND1	1:A:102:LEU:O	2.41	0.45
1:A:620:GLN:HG3	1:A:621:ALA:H	1.81	0.45
1:B:490:VAL:HG11	1:B:533:ILE:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:791:LYS:O	1:A:795:ILE:HG13	2.17	0.44
1:A:627:LEU:HB2	1:A:656:PRO:HA	1.99	0.44
1:B:513:ILE:HG21	1:B:578:ILE:HD13	1.99	0.44
1:A:1226:MET:HB3	1:A:1230:TRP:CZ3	2.52	0.44
1:A:820:THR:HG23	1:A:893:VAL:HG21	1.98	0.44
1:A:1738:TRP:HH2	1:A:1786:PRO:O	2.00	0.44
1:A:1560:GLY:C	1:A:1601:LEU:HD12	2.38	0.43
1:A:1661:ALA:HB2	1:A:1688:PRO:HG2	2.00	0.43
1:B:516:LEU:HB2	1:B:532:PHE:CE1	2.53	0.43
1:A:453:ASN:CG	1:A:456:PHE:HD2	2.21	0.43
1:B:77:LEU:HD12	1:B:99:THR:HG21	2.00	0.43
1:A:1107:GLU:CB	1:A:1754:PHE:CG	3.01	0.43
1:A:406:LEU:HD23	1:A:406:LEU:HA	1.78	0.43
1:A:874:LEU:HD23	1:A:1135:ILE:HG21	2.01	0.43
1:A:1241:VAL:HB	1:A:1431:VAL:HG22	2.00	0.43
1:A:520:LEU:HB3	1:A:556:HIS:HE2	1.84	0.43
1:B:205:PHE:CE1	1:B:403:PHE:HB3	2.54	0.43
1:A:614:THR:HG21	1:A:673:VAL:HG11	2.01	0.42
1:B:53:TYR:CE2	1:B:55:TYR:HB2	2.54	0.42
1:B:401:ASP:O	1:B:405:HIS:ND1	2.52	0.42
1:A:216:THR:HB	1:A:219:HIS:CG	2.55	0.42
1:A:1457:ASN:ND2	1:A:1518:TYR:OH	2.52	0.42
1:A:1587:HIS:CG	1:A:1588:PRO:HD2	2.54	0.42
1:B:73:HIS:HD2	1:B:77:LEU:HD21	1.84	0.42
1:A:77:LEU:HD12	1:A:99:THR:HG21	2.01	0.42
1:A:1235:PRO:O	1:A:1248:THR:OG1	2.25	0.42
1:A:1390:LEU:O	1:A:1394:GLN:HG2	2.20	0.42
1:A:53:TYR:CE2	1:A:55:TYR:HB2	2.54	0.41
1:A:1177:GLN:HB3	1:A:1198:LEU:HB3	2.02	0.41
1:A:1749:ILE:HG23	1:A:1750:ARG:HG3	2.02	0.41
1:B:216:THR:HB	1:B:219:HIS:CG	2.55	0.41
1:A:114:VAL:HG11	1:A:140:LEU:HD11	2.02	0.41
1:A:1467:VAL:HG21	1:A:1497:VAL:HG22	2.03	0.41
1:A:1539:LEU:HB2	1:A:1542:TYR:CD1	2.55	0.41
1:B:213:VAL:HG22	1:B:231:TYR:HB3	2.02	0.41
1:A:437:ILE:HA	1:A:458:VAL:O	2.20	0.41
1:A:1737:ILE:CD1	1:A:1770:LEU:HD23	2.50	0.41
1:A:1385:ALA:HA	1:A:1568:SER:HA	2.02	0.41
1:A:187:LEU:HD22	1:A:524:PRO:HA	2.02	0.41
1:A:651:PRO:HD2	1:A:654:MET:HB2	2.03	0.40
1:A:1724:ALA:HB3	1:A:1749:ILE:HG22	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:ILE:HG21	1:B:115:GLN:HB2	2.03	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	1783/1814 (98%)	1641 (92%)	140 (8%)	2 (0%)	51	85
1	B	583/1814 (32%)	547 (94%)	35 (6%)	1 (0%)	47	81
All	All	2366/3628 (65%)	2188 (92%)	175 (7%)	3 (0%)	51	85

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	488	ILE
1	B	488	ILE
1	A	1517	ILE

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	1527/1562 (98%)	1509 (99%)	18 (1%)	71	84
1	B	499/1562 (32%)	496 (99%)	3 (1%)	86	92

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2026/3124 (65%)	2005 (99%)	21 (1%)	76 86

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	393	PHE
1	A	406	LEU
1	A	572	PHE
1	A	775	LEU
1	A	792	ARG
1	A	873	LEU
1	A	1106	ARG
1	A	1280	ARG
1	A	1314	PHE
1	A	1326	SER
1	A	1405	PHE
1	A	1411	TYR
1	A	1427	ARG
1	A	1458	PHE
1	A	1518	TYR
1	A	1547	ILE
1	A	1764	LEU
1	A	1794	PHE
1	B	254	ASP
1	B	393	PHE
1	B	572	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	HIS
1	A	150	HIS
1	A	177	ASN
1	A	211	GLN
1	A	957	GLN
1	A	1089	GLN
1	A	1177	GLN
1	A	1184	GLN
1	A	1325	GLN
1	A	1359	GLN
1	A	1508	ASN

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Mol	Chain	Res	Type
1	A	1609	ASN
1	A	1618	HIS
1	B	26	HIS
1	B	73	HIS
1	B	150	HIS
1	B	211	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
2	PNS	A	1902	1	0,3,21	-	-	0,3,29	-	-
2	PNS	A	1901	1	0,3,21	-	-	0,3,29	-	-

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.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.