



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 01:28 PM EDT

PDB ID : 8EIL / pdb_00008eil
Title : C-Terminal Domain of BrxL from Acinetobacter BREX type I phage restriction system
Authors : Doyle, L.A.; Stoddard, B.L.; Kaiser, B.
Deposited on : 2022-09-15
Resolution : 2.25 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

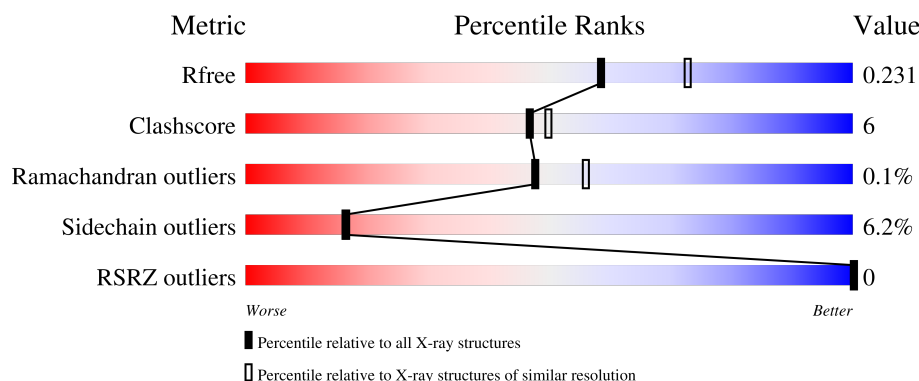
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	186	
1	B	186	
1	C	186	
1	D	186	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SIN	D	701	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5208 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.

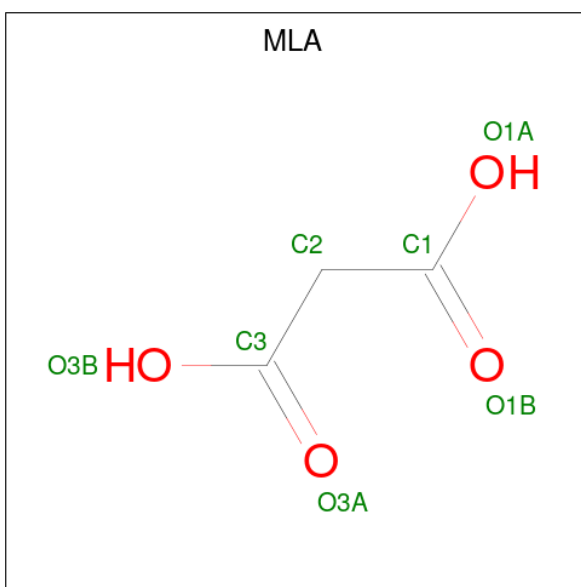
- Molecule 1 is a protein called Protease Lon-related BREX system protein BrxL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	0	0	0
			1290	833	208	243	6			
1	B	180	Total	C	N	O	S	0	0	0
			1300	838	209	246	7			
1	C	178	Total	C	N	O	S	0	0	0
			1243	799	206	233	5			
1	D	177	Total	C	N	O	S	0	0	0
			1233	792	205	231	5			

There are 16 discrepancies between the modelled and reference sequences:

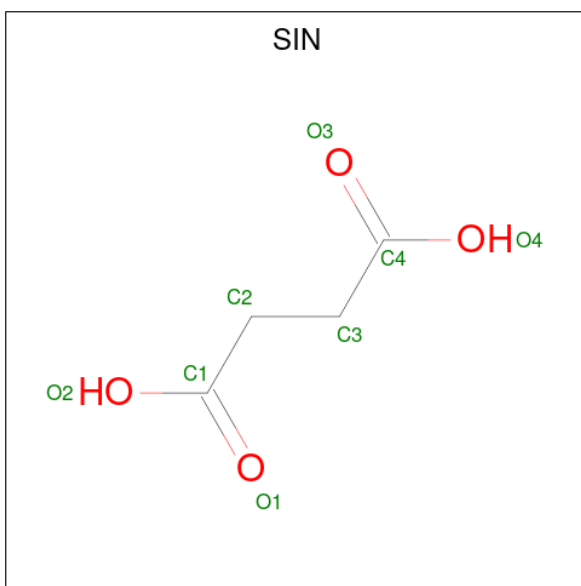
Chain	Residue	Modelled	Actual	Comment	Reference
A	494	GLY	-	expression tag	UNP A0A3R9EDI8
A	495	SER	-	expression tag	UNP A0A3R9EDI8
A	496	HIS	-	expression tag	UNP A0A3R9EDI8
A	497	MET	-	expression tag	UNP A0A3R9EDI8
B	494	GLY	-	expression tag	UNP A0A3R9EDI8
B	495	SER	-	expression tag	UNP A0A3R9EDI8
B	496	HIS	-	expression tag	UNP A0A3R9EDI8
B	497	MET	-	expression tag	UNP A0A3R9EDI8
C	494	GLY	-	expression tag	UNP A0A3R9EDI8
C	495	SER	-	expression tag	UNP A0A3R9EDI8
C	496	HIS	-	expression tag	UNP A0A3R9EDI8
C	497	MET	-	expression tag	UNP A0A3R9EDI8
D	494	GLY	-	expression tag	UNP A0A3R9EDI8
D	495	SER	-	expression tag	UNP A0A3R9EDI8
D	496	HIS	-	expression tag	UNP A0A3R9EDI8
D	497	MET	-	expression tag	UNP A0A3R9EDI8

- Molecule 2 is MALONIC ACID (CCD ID: MLA) (formula: C₃H₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	3	4		
2	B	1	Total	C	O	0	0
			7	3	4		

- Molecule 3 is SUCCINIC ACID (CCD ID: SIN) (formula: $C_4H_6O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			8	4	4		

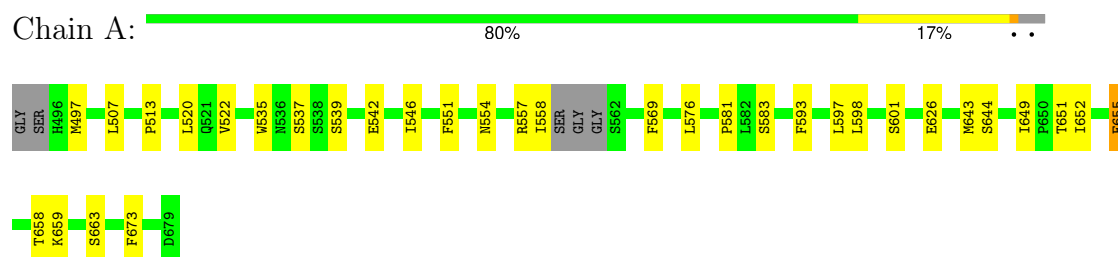
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	40	Total 40	O 40	0	0
4	B	35	Total 35	O 35	0	0
4	C	24	Total 24	O 24	0	0
4	D	21	Total 21	O 21	0	0

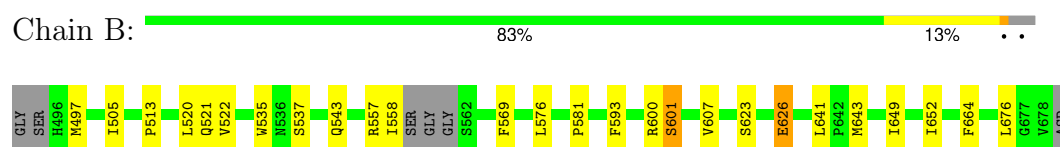
3 Residue-property plots [i](#)

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.

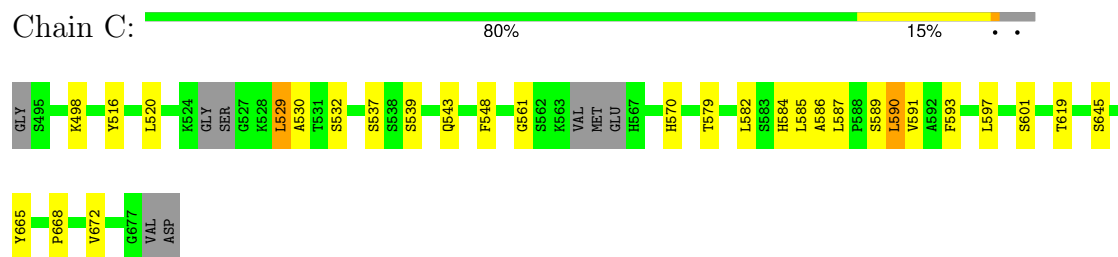
- Molecule 1: Protease Lon-related BREX system protein BrxL



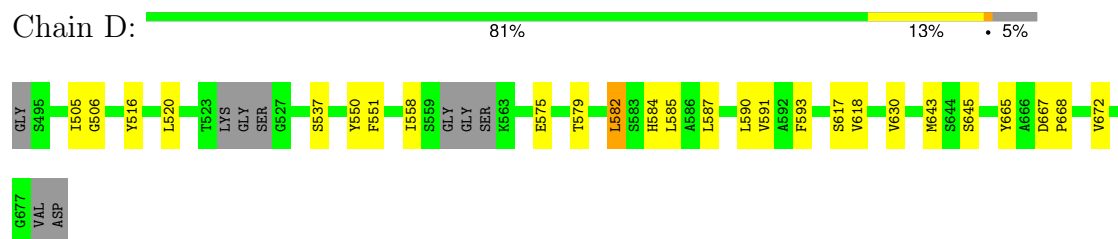
- Molecule 1: Protease Lon-related BREX system protein BrxL



- Molecule 1: Protease Lon-related BREX system protein BrxL



- Molecule 1: Protease Lon-related BREX system protein BrxL



4 Data and refinement statistics

Property	Value	Source
Space group	F 2 3	Depositor
Cell constants a, b, c, α , β , γ	213.34Å 213.34Å 213.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.94 – 2.25 48.94 – 2.25	Depositor EDS
% Data completeness (in resolution range)	92.6 (48.94-2.25) 92.1 (48.94-2.25)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.18_3855	Depositor
R, R_{free}	0.216 , 0.230 0.216 , 0.231	Depositor DCC
R_{free} test set	2024 reflections (5.33%)	wwPDB-VP
Wilson B-factor (Å ²)	45.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 63.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.499 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5208	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.16	0/1317	0.34	0/1794
1	B	0.16	0/1327	0.37	0/1807
1	C	0.16	0/1270	0.39	0/1731
1	D	0.15	0/1260	0.33	0/1719
All	All	0.16	0/5174	0.36	0/7051

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 [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	A	1290	0	1226	16	0
1	B	1300	0	1252	13	0
1	C	1243	0	1144	14	0
1	D	1233	0	1119	15	0
2	A	7	0	2	0	0
2	B	7	0	2	0	0
3	D	8	0	4	0	0
4	A	40	0	0	0	0
4	B	35	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	24	0	0	0	0
4	D	21	0	0	0	0
All	All	5208	0	4749	56	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 6.

All (56) 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:D:520:LEU:HD22	1:D:593:PHE:HA	1.67	0.74
1:C:520:LEU:HD22	1:C:593:PHE:HA	1.70	0.73
1:D:516:TYR:HD1	1:D:575:GLU:HA	1.59	0.67
1:C:516:TYR:OH	1:C:579:THR:OG1	2.14	0.65
1:D:582:LEU:HD22	1:D:584:HIS:H	1.62	0.65
1:A:497:MET:O	1:A:601:SER:OG	2.15	0.64
1:D:516:TYR:OH	1:D:579:THR:OG1	2.18	0.60
1:A:520:LEU:HD11	1:A:569:PHE:HB3	1.87	0.57
1:B:600:ARG:HB2	1:B:676:LEU:HD22	1.86	0.55
1:B:505:ILE:HD12	1:B:607:VAL:HG11	1.88	0.55
1:A:520:LEU:HD22	1:A:593:PHE:HA	1.90	0.52
1:C:529:LEU:HD23	1:C:548:PHE:CG	2.45	0.52
1:C:585:LEU:O	1:C:589:SER:OG	2.22	0.52
1:B:649:ILE:HA	1:B:652:ILE:HD12	1.92	0.51
1:C:582:LEU:HD23	1:C:585:LEU:HG	1.92	0.51
1:A:649:ILE:HA	1:A:652:ILE:HD12	1.93	0.50
1:B:520:LEU:HD11	1:B:569:PHE:HB3	1.93	0.50
1:C:587:LEU:HD21	1:C:665:TYR:CE1	2.46	0.50
1:B:557:ARG:O	1:B:558:ILE:HG13	2.11	0.50
1:D:506:GLY:HA3	1:D:582:LEU:HD21	1.95	0.49
1:D:587:LEU:HD21	1:D:665:TYR:CE2	2.49	0.48
1:D:591:VAL:HA	1:D:672:VAL:HG22	1.96	0.48
1:A:554:ASN:HB3	1:A:557:ARG:HD3	1.94	0.48
1:B:513:PRO:HG3	1:B:626:GLU:HB2	1.96	0.48
1:B:497:MET:O	1:B:601:SER:HB3	2.13	0.48
1:C:543:GLN:NE2	1:C:586:ALA:H	2.12	0.48
1:C:520:LEU:HD13	1:C:593:PHE:CD1	2.50	0.47
1:C:530:ALA:HB3	1:C:570:HIS:HD2	1.80	0.47
1:D:516:TYR:CE2	1:D:582:LEU:HG	2.50	0.47
1:C:590:LEU:HD22	1:C:668:PRO:HB3	1.96	0.46
1:A:542:GLU:O	1:A:546:ILE:HG13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:516:TYR:CD1	1:D:575:GLU:HA	2.46	0.45
1:A:535:TRP:CE2	1:A:581:PRO:HA	2.52	0.45
1:A:513:PRO:HG3	1:A:626:GLU:HB3	1.99	0.44
1:A:658:THR:HG21	1:D:643:MET:HG3	1.99	0.44
1:C:520:LEU:HD13	1:C:593:PHE:CE1	2.52	0.44
1:D:550:TYR:CZ	1:D:668:PRO:HG2	2.53	0.44
1:A:558:ILE:HG22	1:A:598:LEU:HD11	2.00	0.44
1:B:520:LEU:HD22	1:B:593:PHE:HA	2.00	0.44
1:A:576:LEU:HD23	1:A:576:LEU:HA	1.86	0.44
1:B:535:TRP:CE2	1:B:581:PRO:HA	2.53	0.43
1:A:507:LEU:O	1:A:583:SER:OG	2.35	0.43
1:A:551:PHE:CZ	1:A:597:LEU:HD12	2.53	0.43
1:D:520:LEU:HD13	1:D:593:PHE:CE1	2.52	0.43
1:B:641:LEU:O	1:B:664:PHE:HA	2.18	0.43
1:B:543:GLN:NE2	4:B:801:HOH:O	2.42	0.43
1:B:576:LEU:HD23	1:B:576:LEU:HA	1.90	0.43
1:C:591:VAL:HA	1:C:672:VAL:HG22	1.99	0.43
1:D:505:ILE:HD13	1:D:630:VAL:HG12	2.00	0.43
1:A:655:GLU:O	1:A:659:LYS:HG2	2.19	0.42
1:A:522:VAL:HG22	1:A:569:PHE:CD1	2.54	0.42
1:A:673:PHE:CD2	1:C:498:LYS:HE2	2.54	0.42
1:B:522:VAL:HG22	1:B:569:PHE:CD1	2.55	0.41
1:C:582:LEU:HG	1:C:584:HIS:H	1.86	0.41
1:D:551:PHE:CD2	1:D:593:PHE:HB3	2.56	0.41
1:D:618:VAL:HG13	1:D:668:PRO:HD3	2.04	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	177/186 (95%)	176 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	176/186 (95%)	173 (98%)	3 (2%)	0	100	100
1	C	172/186 (92%)	167 (97%)	4 (2%)	1 (1%)	21	21
1	D	171/186 (92%)	168 (98%)	3 (2%)	0	100	100
All	All	696/744 (94%)	684 (98%)	11 (2%)	1 (0%)	48	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	561	GLY

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	126/150 (84%)	119 (94%)	7 (6%)	19	20
1	B	132/150 (88%)	126 (96%)	6 (4%)	24	29
1	C	116/150 (77%)	107 (92%)	9 (8%)	11	10
1	D	113/150 (75%)	105 (93%)	8 (7%)	13	12
All	All	487/600 (81%)	457 (94%)	30 (6%)	16	16

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	537	SER
1	A	539	SER
1	A	643	MET
1	A	644	SER
1	A	651	THR
1	A	655	GLU
1	A	663	SER
1	B	521	GLN
1	B	537	SER
1	B	601	SER

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Mol	Chain	Res	Type
1	B	623	SER
1	B	626	GLU
1	B	643	MET
1	C	529	LEU
1	C	532	SER
1	C	537	SER
1	C	539	SER
1	C	590	LEU
1	C	597	LEU
1	C	601	SER
1	C	619	THR
1	C	645	SER
1	D	537	SER
1	D	558	ILE
1	D	582	LEU
1	D	585	LEU
1	D	590	LEU
1	D	617	SER
1	D	645	SER
1	D	667	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	577	GLN
1	A	603	GLN
1	B	603	GLN
1	B	661	GLN
1	C	543	GLN
1	C	570	HIS
1	C	605	GLN
1	D	543	GLN
1	D	554	ASN

5.3.3 RNA ⓘ

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](#)

3 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	MLA	A	701	-	6,6,6	1.37	0	7,7,7	1.30	0
3	SIN	D	701	-	7,7,7	1.04	0	8,8,8	1.66	3 (37%)
2	MLA	B	701	-	6,6,6	1.42	0	7,7,7	1.14	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
2	MLA	A	701	-	-	2/4/4/4	-
3	SIN	D	701	-	-	5/5/5/5	-
2	MLA	B	701	-	-	2/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	701	SIN	C2-C3-C4	-2.04	108.25	113.67
3	D	701	SIN	O2-C1-C2	2.03	120.42	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	701	SIN	O4-C4-C3	2.02	120.39	114.00

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	701	SIN	C1-C2-C3-C4
2	B	701	MLA	O1A-C1-C2-C3
2	B	701	MLA	O1B-C1-C2-C3
2	A	701	MLA	O1B-C1-C2-C3
2	A	701	MLA	O1A-C1-C2-C3
3	D	701	SIN	C2-C3-C4-O4
3	D	701	SIN	C2-C3-C4-O3
3	D	701	SIN	O2-C1-C2-C3
3	D	701	SIN	O1-C1-C2-C3

There are no ring outliers.

No monomer is involved in short contacts.

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	181/186 (97%)	-1.21	0 100 100	32, 53, 82, 98	0
1	B	180/186 (96%)	-1.20	0 100 100	34, 53, 84, 95	0
1	C	178/186 (95%)	-1.03	0 100 100	34, 71, 101, 115	0
1	D	177/186 (95%)	-1.05	0 100 100	35, 71, 98, 121	0
All	All	716/744 (96%)	-1.12	0 100 100	32, 61, 95, 121	0

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
2	MLA	A	701	7/7	0.98	0.03	57,62,73,74	0
2	MLA	B	701	7/7	0.98	0.04	59,61,71,72	0
3	SIN	D	701	8/8	0.99	0.03	55,67,68,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.