



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:03 AM UTC

PDB ID : 4FN5 / pdb_00004fn5
Title : ELONGATION FACTOR G 1 (PSEUDOMONAS AERUGINOSA) IN COM-
PLEX WITH Argyrin B
Authors : Kirby, C.A.; Palestrant, D.
Deposited on : 2012-06-19
Resolution : 2.90 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)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

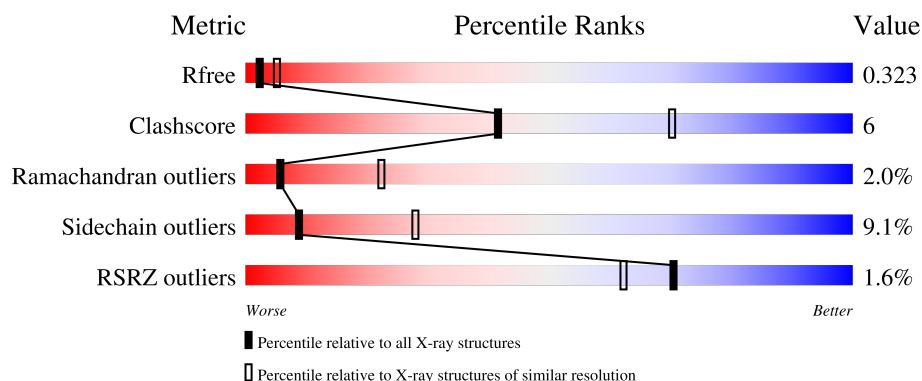
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

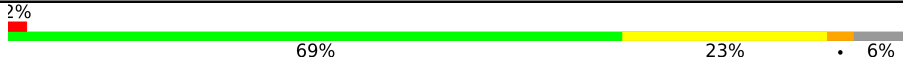
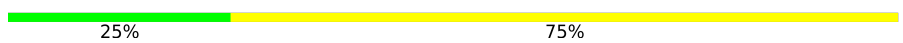
The reported resolution of this entry is 2.90 Å.

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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	709	
2	B	8	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5253 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 Elongation factor G 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	670	Total	C	N	O	S	0	0	0
			5177	3252	893	1005	27			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q9HWD2
A	-1	SER	-	expression tag	UNP Q9HWD2
A	0	HIS	-	expression tag	UNP Q9HWD2

- Molecule 2 is a protein called Argyrin B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	8	Total	C	N	O	S	0	0	0
			60	41	10	8	1			

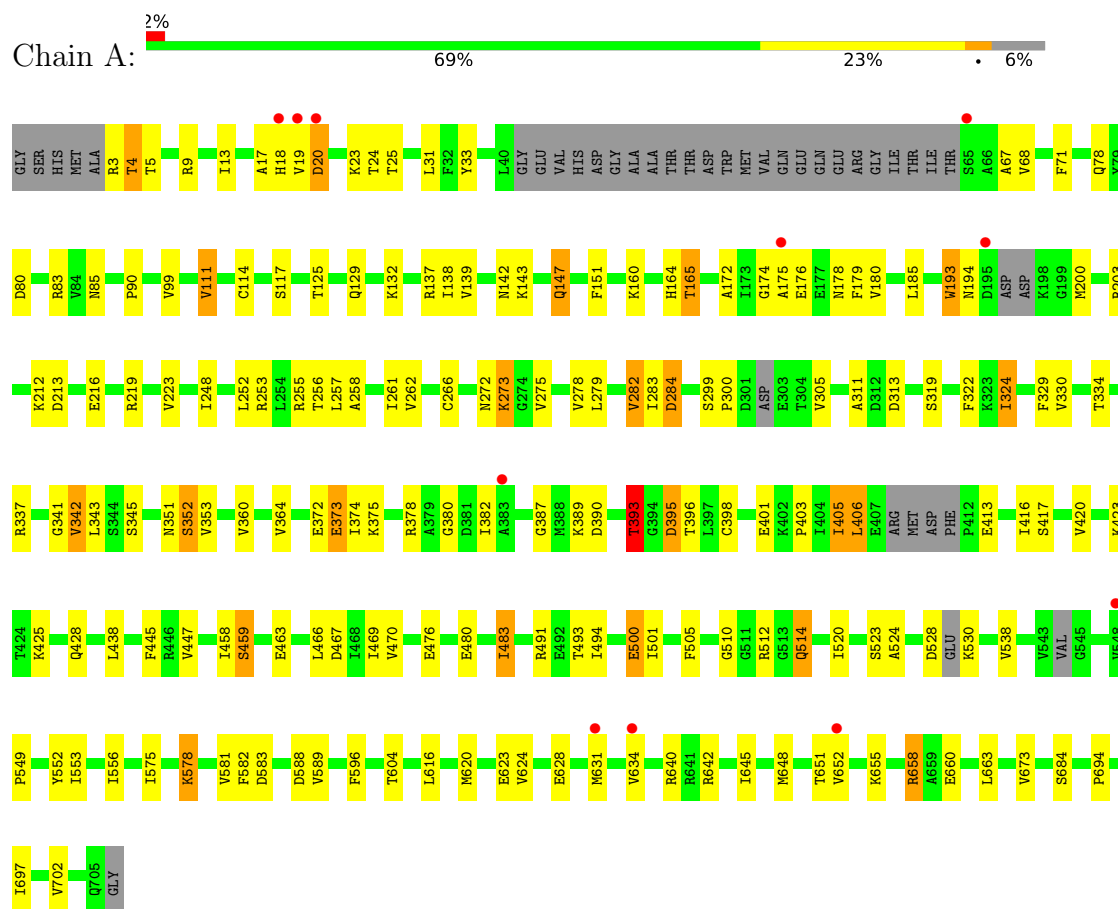
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	16	Total	O	0	0
			16	16		

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: Elongation factor G 1



• Molecule 2: Argirin B



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	125.78Å 88.10Å 74.32Å 90.00° 107.91° 90.00°	Depositor
Resolution (Å)	55.39 – 2.90 55.39 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.1 (55.39-2.90) 98.5 (55.39-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.91Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.202 , 0.300 0.214 , 0.323	Depositor DCC
R_{free} test set	857 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	67.5	Xtriage
Anisotropy	0.316	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 64.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5253	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% 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: 0UO, DHA, BB9, DBB, DAL, SAR

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	0.90	0/5257	1.48	38/7088 (0.5%)
2	B	0.77	0/18	1.28	0/22
All	All	0.90	0/5275	1.48	38/7110 (0.5%)

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	463	GLU	N-CA-C	-7.90	102.61	111.07
1	A	20	ASP	CA-CB-CG	6.99	119.59	112.60
1	A	111	VAL	CB-CA-C	6.82	119.91	111.25
1	A	549	PRO	CA-C-N	6.81	129.30	120.44
1	A	549	PRO	C-N-CA	6.81	129.30	120.44
1	A	628	GLU	N-CA-C	6.75	119.22	111.11
1	A	262	VAL	N-CA-CB	6.74	118.30	111.36
1	A	642	ARG	CA-C-N	6.38	126.16	120.10
1	A	642	ARG	C-N-CA	6.38	126.16	120.10
1	A	160	LYS	N-CA-C	6.34	119.17	111.82
1	A	18	HIS	N-CA-C	6.17	120.54	111.34
1	A	179	PHE	N-CA-C	6.15	118.47	108.63
1	A	645	ILE	N-CA-C	5.90	116.63	108.84
1	A	552	TYR	CA-C-N	5.89	126.09	120.43
1	A	552	TYR	C-N-CA	5.89	126.09	120.43
1	A	193	TRP	N-CA-C	5.77	118.96	109.72
1	A	329	PHE	CA-C-N	5.69	128.26	120.46
1	A	329	PHE	C-N-CA	5.69	128.26	120.46
1	A	640	ARG	N-CA-C	-5.56	106.47	113.20
1	A	395	ASP	CA-CB-CG	5.52	118.12	112.60
1	A	352	SER	CA-C-N	5.47	127.54	120.70
1	A	352	SER	C-N-CA	5.47	127.54	120.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	284	ASP	N-CA-C	5.47	117.05	111.14
1	A	78	GLN	N-CA-C	5.43	119.16	112.54
1	A	172	ALA	N-CA-C	5.39	118.60	109.76
1	A	596	PHE	CA-CB-CG	-5.39	108.41	113.80
1	A	282	VAL	N-CA-C	-5.33	105.30	110.42
1	A	398	CYS	N-CA-C	5.23	116.61	109.18
1	A	256	THR	N-CA-C	-5.22	105.50	111.14
1	A	393	THR	CB-CA-C	5.11	118.39	109.65
1	A	273	LYS	N-CA-C	5.10	117.61	108.17
1	A	165	THR	CB-CA-C	5.06	118.97	110.97
1	A	342	VAL	N-CA-CB	5.06	120.19	111.39
1	A	278	VAL	N-CA-C	-5.04	105.79	110.53
1	A	467	ASP	CA-C-N	5.02	126.89	120.56
1	A	467	ASP	C-N-CA	5.02	126.89	120.56
1	A	334	THR	CB-CA-C	5.01	117.93	110.62
1	A	648	MET	N-CA-C	5.01	116.21	107.49

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	5177	0	5179	66	0
2	B	60	0	43	1	0
3	A	16	0	0	0	0
All	All	5253	0	5222	67	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 (67) 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:3:ARG:HG2	1:A:4:THR:H	1.43	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ALA:HB3	1:A:23:LYS:HB2	1.69	0.75
1:A:343:LEU:HD11	1:A:360:VAL:HG21	1.76	0.66
1:A:194:ASN:HD21	1:A:203:ARG:HG3	1.64	0.62
1:A:364:VAL:HG12	1:A:373:GLU:HA	1.81	0.62
1:A:151:PHE:CD1	1:A:266:CYS:HB3	2.35	0.61
1:A:423:LYS:HG3	1:A:480:GLU:HG2	1.84	0.59
1:A:396:THR:HG23	1:A:406:LEU:HD22	1.85	0.59
1:A:20:ASP:O	1:A:114:CYS:HB2	2.04	0.58
1:A:300:PRO:HG2	1:A:403:PRO:HB2	1.86	0.57
1:A:493:THR:HB	1:A:616:LEU:HD11	1.86	0.57
1:A:193:TRP:HB2	1:A:273:LYS:HG2	1.88	0.55
1:A:85:ASN:HB3	1:A:382:ILE:HD12	1.88	0.55
1:A:114:CYS:HB3	1:A:117:SER:OG	2.09	0.53
1:A:337:ARG:HH11	1:A:380:GLY:HA2	1.73	0.52
1:A:255:ARG:HB2	1:A:261:ILE:HG12	1.90	0.52
1:A:631:MET:HG3	1:A:655:LYS:HG3	1.92	0.51
1:A:253:ARG:HD2	1:A:284:ASP:O	2.12	0.50
1:A:174:GLY:HA3	1:A:178:ASN:O	2.12	0.49
1:A:142:ASN:HD22	1:A:143:LYS:HG3	1.78	0.49
1:A:624:VAL:HG11	1:A:634:VAL:HG11	1.94	0.49
1:A:279:LEU:O	1:A:282:VAL:HB	2.13	0.48
1:A:505:PHE:CE1	1:A:514:GLN:HB3	2.49	0.47
1:A:420:VAL:HG23	1:A:483:ILE:HG22	1.95	0.47
1:A:520:ILE:HA	1:A:581:VAL:HA	1.96	0.47
1:A:620:MET:SD	1:A:663:LEU:HD13	2.55	0.47
1:A:31:LEU:CD1	1:A:68:VAL:HG11	2.45	0.47
1:A:447:VAL:HG22	1:A:458:ILE:HG22	1.97	0.47
1:A:353:VAL:HG21	1:A:395:ASP:HA	1.97	0.47
1:A:253:ARG:O	1:A:257:LEU:HG	2.16	0.46
1:A:313:ASP:O	1:A:341:GLY:HA2	2.16	0.46
1:A:417:SER:OG	1:A:459:SER:HB3	2.16	0.46
1:A:623:GLU:HG3	1:A:658:ARG:HG3	1.97	0.45
1:A:694:PRO:HG2	1:A:697:ILE:HG12	1.97	0.45
1:A:71:PHE:CE2	1:A:83:ARG:HG2	2.52	0.45
1:A:5:THR:HG21	1:A:83:ARG:HG3	1.98	0.45
1:A:494:ILE:HG13	1:A:524:ALA:HA	1.99	0.45
1:A:147:GLN:HA	1:A:176:GLU:HG2	1.99	0.45
1:A:212:LYS:O	1:A:216:GLU:HG2	2.17	0.45
1:A:500:GLU:HB3	1:A:582:PHE:HZ	1.83	0.44
1:A:553:ILE:HD13	1:A:556:ILE:HD12	1.98	0.44
1:A:500:GLU:HB3	1:A:582:PHE:CZ	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:ALA:HB2	1:A:322:PHE:HZ	1.82	0.44
1:A:319:SER:HB2	1:A:337:ARG:O	2.18	0.44
1:A:90:PRO:HD2	1:A:99:VAL:HG23	1.99	0.44
1:A:3:ARG:CG	1:A:4:THR:H	2.22	0.43
1:A:470:VAL:HG21	1:A:483:ILE:HG21	2.00	0.43
1:A:458:ILE:HD12	1:A:466:LEU:HD22	2.00	0.43
1:A:438:LEU:HD13	1:A:469:ILE:HG23	2.00	0.43
2:B:1:DAL:HB3	2:B:6:DBB:HB3	2.01	0.42
1:A:255:ARG:HA	1:A:258:ALA:HB3	2.00	0.42
1:A:324:ILE:HD11	1:A:393:THR:N	2.35	0.42
1:A:3:ARG:HG3	1:A:378:ARG:HD2	2.03	0.41
1:A:33:TYR:CD1	1:A:200:MET:HE2	2.56	0.41
1:A:31:LEU:HD13	1:A:68:VAL:HG11	2.02	0.41
1:A:9:ARG:NE	1:A:80:ASP:HB2	2.36	0.41
1:A:9:ARG:HE	1:A:80:ASP:HB2	1.86	0.41
1:A:71:PHE:HE2	1:A:83:ARG:HG2	1.85	0.41
1:A:528:ASP:O	1:A:530:LYS:HG3	2.20	0.41
1:A:129:GLN:O	1:A:132:LYS:HB3	2.21	0.40
1:A:523:SER:HB2	1:A:578:LYS:HB2	2.03	0.40
1:A:445:PHE:CE1	1:A:458:ILE:HB	2.56	0.40
1:A:13:ILE:HG21	1:A:282:VAL:HG13	2.03	0.40
1:A:491:ARG:HD3	1:A:575:ILE:HG12	2.04	0.40
1:A:219:ARG:O	1:A:223:VAL:HG23	2.21	0.40
1:A:248:ILE:O	1:A:252:LEU:HG	2.21	0.40
1:A:300:PRO:HG3	1:A:405:ILE:HG13	2.03	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	656/709 (92%)	599 (91%)	44 (7%)	13 (2%)	6	22

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	2/8 (25%)	2 (100%)	0	0	100	100
All	All	658/717 (92%)	601 (91%)	44 (7%)	13 (2%)	6	22

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	352	SER
1	A	372	GLU
1	A	389	LYS
1	A	428	GLN
1	A	588	ASP
1	A	652	VAL
1	A	311	ALA
1	A	387	GLY
1	A	425	LYS
1	A	476	GLU
1	A	272	ASN
1	A	175	ALA
1	A	510	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	558/589 (95%)	507 (91%)	51 (9%)	9	28
2	B	1/1 (100%)	1 (100%)	0	100	100
All	All	559/590 (95%)	508 (91%)	51 (9%)	9	28

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	19	VAL
1	A	24	THR

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Mol	Chain	Res	Type
1	A	25	THR
1	A	111	VAL
1	A	125	THR
1	A	137	ARG
1	A	138	ILE
1	A	139	VAL
1	A	147	GLN
1	A	164	HIS
1	A	165	THR
1	A	180	VAL
1	A	185	LEU
1	A	213	ASP
1	A	275	VAL
1	A	283	ILE
1	A	299	SER
1	A	305	VAL
1	A	324	ILE
1	A	330	VAL
1	A	342	VAL
1	A	345	SER
1	A	351	ASN
1	A	373	GLU
1	A	374	ILE
1	A	375	LYS
1	A	390	ASP
1	A	393	THR
1	A	401	GLU
1	A	405	ILE
1	A	406	LEU
1	A	413	GLU
1	A	416	ILE
1	A	459	SER
1	A	483	ILE
1	A	500	GLU
1	A	501	ILE
1	A	512	ARG
1	A	514	GLN
1	A	538	VAL
1	A	578	LYS
1	A	583	ASP
1	A	589	VAL
1	A	604	THR

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Mol	Chain	Res	Type
1	A	651	THR
1	A	658	ARG
1	A	660	GLU
1	A	673	VAL
1	A	684	SER
1	A	702	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	92	HIS
1	A	122	GLN
1	A	369	ASN
1	A	440	GLN
1	A	541	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 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	BB9	B	2	2	2,5,6	1.89	1 (50%)	1,5,7	1.06	0
2	DHA	B	7	2	3,4,5	1.05	0	2,4,6	3.27	2 (100%)
2	0UO	B	4	2	16,17,18	1.19	3 (18%)	17,23,25	1.30	2 (11%)
2	SAR	B	8	2	3,4,5	0.93	0	1,3,5	2.63	1 (100%)
2	DBB	B	6	2	4,5,6	0.55	0	1,5,7	0.76	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	BB9	B	2	2	-	0/0/4/6	-
2	DHA	B	7	2	-	0/0/2/4	-
2	0UO	B	4	2	-	0/7/8/10	0/2/2/2
2	SAR	B	8	2	-	1/1/2/3	-
2	DBB	B	6	2	-	2/3/4/6	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	BB9	C-CA	2.45	1.49	1.45
2	B	4	0UO	C18-C17	2.13	1.44	1.40
2	B	4	0UO	C16-C17	2.13	1.44	1.39
2	B	4	0UO	C18-C13	2.03	1.44	1.41

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	7	DHA	CB-CA-N	-3.51	117.31	125.76
2	B	7	DHA	O-C-CA	-3.02	119.87	125.53
2	B	4	0UO	C9-C10-C11	-2.63	121.83	126.97
2	B	8	SAR	O-C-CA	-2.63	116.13	125.17
2	B	4	0UO	C13-N6-C11	2.30	111.12	109.08

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	8	SAR	C-CA-N-CN
2	B	6	DBB	C-CA-CB-CG
2	B	6	DBB	N-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	6	DBB	1	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	670/709 (94%)	0.49	11 (1%) 70 62	43, 76, 105, 121	0
2	B	2/8 (25%)	-0.46	0 100 100	43, 43, 43, 52	0
All	All	672/717 (93%)	0.49	11 (1%) 70 62	43, 76, 105, 121	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	383	ALA	2.9
1	A	20	ASP	2.7
1	A	652	VAL	2.6
1	A	18	HIS	2.5
1	A	195	ASP	2.4
1	A	65	SER	2.3
1	A	548	VAL	2.3
1	A	19	VAL	2.2
1	A	634	VAL	2.2
1	A	631	MET	2.1
1	A	175	ALA	2.0

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	SAR	B	8	5/6	0.95	0.11	49,49,54,54	0
2	OUO	B	4	16/17	0.96	0.08	44,48,50,50	0
2	DBB	B	6	6/7	0.96	0.09	48,50,53,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BB9	B	2	6/7	0.97	0.07	45,45,46,47	0
2	DAL	B	1	4/6	0.97	0.06	36,41,43,45	0
2	DHA	B	7	5/6	0.97	0.09	52,54,57,60	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.