



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:45 PM UTC

PDB ID : 6G1F / pdb_00006g1f
Title : Crystal structure of D-phenylglycine aminotransferase (D-PhgAT) from *Pseudomonas stutzeri* with PLP internal aldimine
Authors : Serpico, A.; Marles-Wright, J.; Campopiano, D.J.
Deposited on : 2018-03-21
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

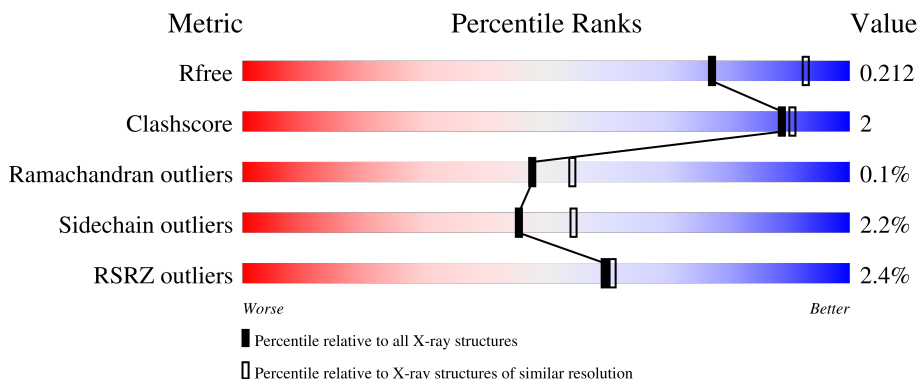
1 Overall quality at a glance

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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	464	2% 90% 5% 5%
1	B	464	2% 88% 7% 5%
1	C	464	2% 88% 7% 5%
1	D	464	0% 90% 5% 5%
1	E	464	3% 88% 5% 6%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	464	5% 87% 7% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 40487 atoms, of which 19845 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 D-phenylglycine aminotransferase.

Mol	Chain	Residues	Atoms								ZeroOcc	AltConf	Trace
1	A	443	Total	C	H	N	O	P	S		0	0	0
			6726	2141	3329	618	625	1	12				
1	B	441	Total	C	H	N	O	P	S		0	0	0
			6690	2127	3313	615	622	1	12				
1	C	440	Total	C	H	N	O	P	S		0	0	0
			6683	2125	3310	614	621	1	12				
1	D	441	Total	C	H	N	O	P	S		0	0	0
			6689	2127	3312	615	622	1	12				
1	E	438	Total	C	H	N	O	P	S		0	0	0
			6653	2116	3297	611	616	1	12				
1	F	438	Total	C	H	N	O	P	S		0	0	0
			6637	2111	3284	611	618	1	12				

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	initiating methionine	UNP Q6VY99
A	-9	SER	-	expression tag	UNP Q6VY99
A	-8	HIS	-	expression tag	UNP Q6VY99
A	-7	HIS	-	expression tag	UNP Q6VY99
A	-6	HIS	-	expression tag	UNP Q6VY99
A	-5	HIS	-	expression tag	UNP Q6VY99
A	-4	HIS	-	expression tag	UNP Q6VY99
A	-3	HIS	-	expression tag	UNP Q6VY99
A	-2	GLY	-	expression tag	UNP Q6VY99
A	-1	SER	-	expression tag	UNP Q6VY99
A	0	SER	-	expression tag	UNP Q6VY99
A	1	GLY	-	expression tag	UNP Q6VY99
B	-10	MET	-	initiating methionine	UNP Q6VY99
B	-9	SER	-	expression tag	UNP Q6VY99
B	-8	HIS	-	expression tag	UNP Q6VY99
B	-7	HIS	-	expression tag	UNP Q6VY99
B	-6	HIS	-	expression tag	UNP Q6VY99

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	HIS	-	expression tag	UNP Q6VY99
B	-4	HIS	-	expression tag	UNP Q6VY99
B	-3	HIS	-	expression tag	UNP Q6VY99
B	-2	GLY	-	expression tag	UNP Q6VY99
B	-1	SER	-	expression tag	UNP Q6VY99
B	0	SER	-	expression tag	UNP Q6VY99
B	1	GLY	-	expression tag	UNP Q6VY99
C	-10	MET	-	initiating methionine	UNP Q6VY99
C	-9	SER	-	expression tag	UNP Q6VY99
C	-8	HIS	-	expression tag	UNP Q6VY99
C	-7	HIS	-	expression tag	UNP Q6VY99
C	-6	HIS	-	expression tag	UNP Q6VY99
C	-5	HIS	-	expression tag	UNP Q6VY99
C	-4	HIS	-	expression tag	UNP Q6VY99
C	-3	HIS	-	expression tag	UNP Q6VY99
C	-2	GLY	-	expression tag	UNP Q6VY99
C	-1	SER	-	expression tag	UNP Q6VY99
C	0	SER	-	expression tag	UNP Q6VY99
C	1	GLY	-	expression tag	UNP Q6VY99
D	-10	MET	-	initiating methionine	UNP Q6VY99
D	-9	SER	-	expression tag	UNP Q6VY99
D	-8	HIS	-	expression tag	UNP Q6VY99
D	-7	HIS	-	expression tag	UNP Q6VY99
D	-6	HIS	-	expression tag	UNP Q6VY99
D	-5	HIS	-	expression tag	UNP Q6VY99
D	-4	HIS	-	expression tag	UNP Q6VY99
D	-3	HIS	-	expression tag	UNP Q6VY99
D	-2	GLY	-	expression tag	UNP Q6VY99
D	-1	SER	-	expression tag	UNP Q6VY99
D	0	SER	-	expression tag	UNP Q6VY99
D	1	GLY	-	expression tag	UNP Q6VY99
E	-10	MET	-	initiating methionine	UNP Q6VY99
E	-9	SER	-	expression tag	UNP Q6VY99
E	-8	HIS	-	expression tag	UNP Q6VY99
E	-7	HIS	-	expression tag	UNP Q6VY99
E	-6	HIS	-	expression tag	UNP Q6VY99
E	-5	HIS	-	expression tag	UNP Q6VY99
E	-4	HIS	-	expression tag	UNP Q6VY99
E	-3	HIS	-	expression tag	UNP Q6VY99
E	-2	GLY	-	expression tag	UNP Q6VY99
E	-1	SER	-	expression tag	UNP Q6VY99
E	0	SER	-	expression tag	UNP Q6VY99

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	1	GLY	-	expression tag	UNP Q6VY99
F	-10	MET	-	initiating methionine	UNP Q6VY99
F	-9	SER	-	expression tag	UNP Q6VY99
F	-8	HIS	-	expression tag	UNP Q6VY99
F	-7	HIS	-	expression tag	UNP Q6VY99
F	-6	HIS	-	expression tag	UNP Q6VY99
F	-5	HIS	-	expression tag	UNP Q6VY99
F	-4	HIS	-	expression tag	UNP Q6VY99
F	-3	HIS	-	expression tag	UNP Q6VY99
F	-2	GLY	-	expression tag	UNP Q6VY99
F	-1	SER	-	expression tag	UNP Q6VY99
F	0	SER	-	expression tag	UNP Q6VY99
F	1	GLY	-	expression tag	UNP Q6VY99

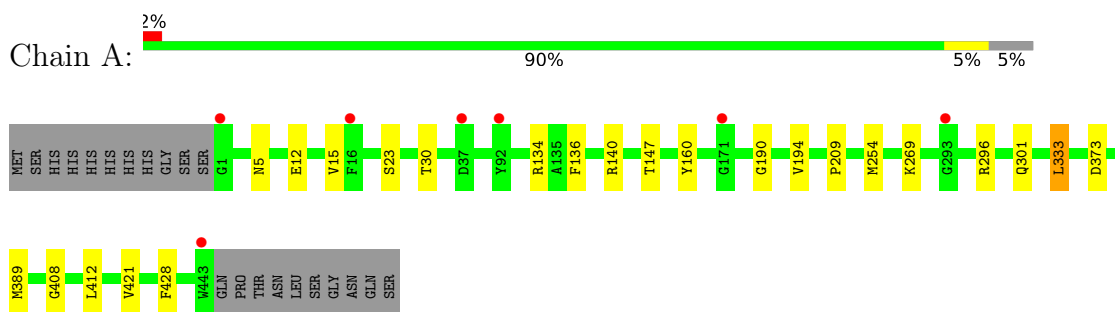
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	106	Total	O	0	0
			106	106		
2	B	120	Total	O	0	0
			120	120		
2	C	59	Total	O	0	0
			59	59		
2	D	83	Total	O	0	0
			83	83		
2	E	35	Total	O	0	0
			35	35		
2	F	6	Total	O	0	0
			6	6		

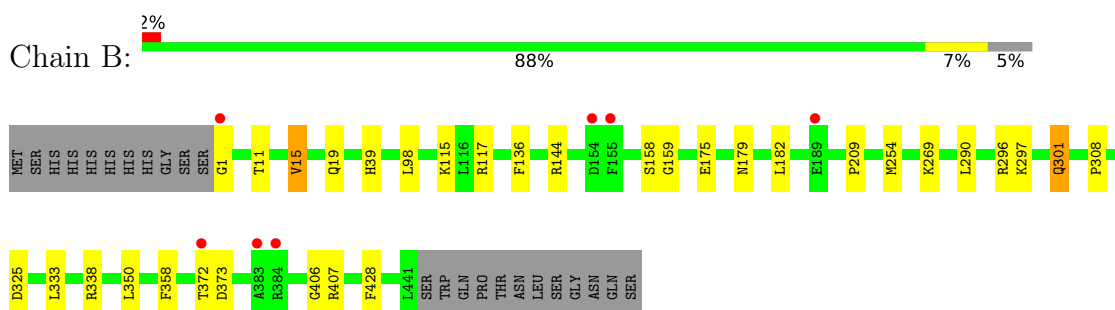
3 Residue-property plots [i](#)

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.

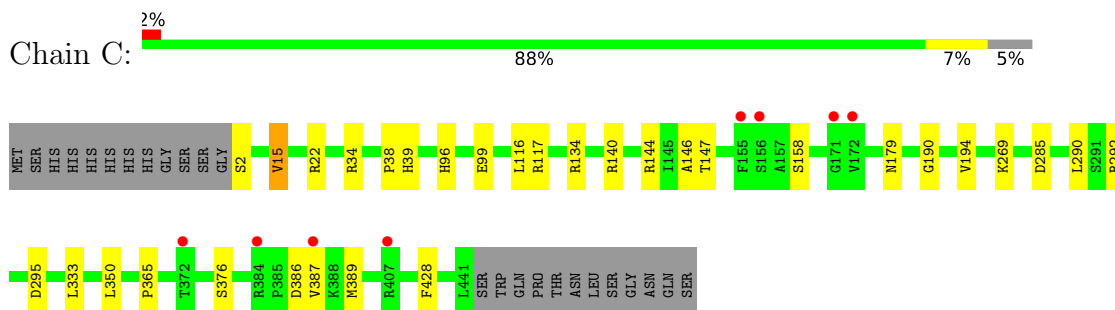
- Molecule 1: D-phenylglycine aminotransferase



- Molecule 1: D-phenylglycine aminotransferase

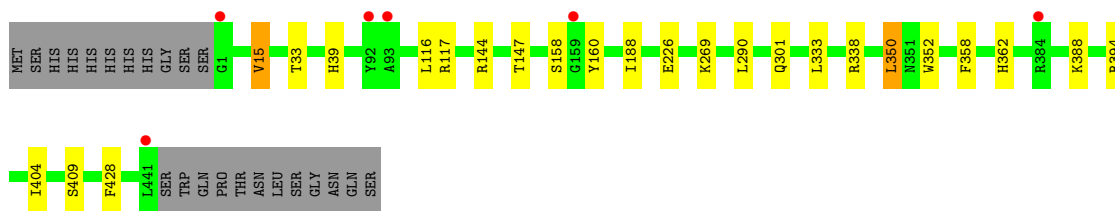


- Molecule 1: D-phenylglycine aminotransferase

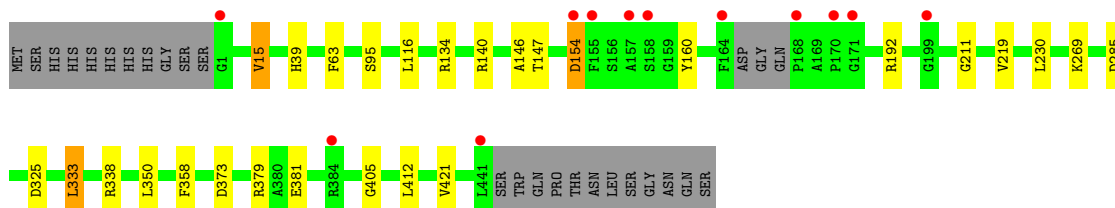
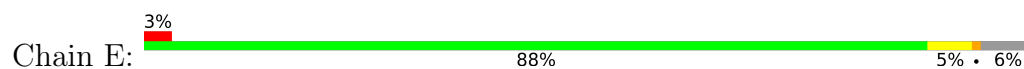


- Molecule 1: D-phenylglycine aminotransferase

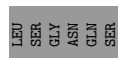
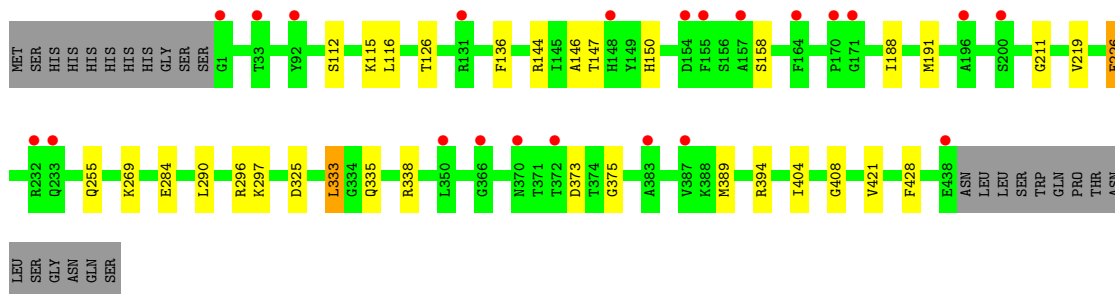
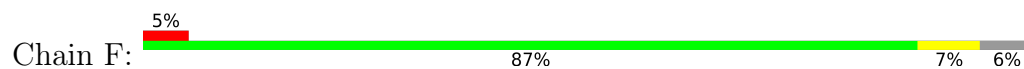




● Molecule 1: D-phenylglycine aminotransferase



● Molecule 1: D-phenylglycine aminotransferase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	329.28Å 83.90Å 133.42Å 90.00° 111.58° 90.00°	Depositor
Resolution (Å)	49.47 – 2.25 49.47 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.47-2.25) 99.8 (49.47-2.25)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 2.25Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.183 , 0.211 0.186 , 0.212	Depositor DCC
R_{free} test set	8147 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	47.8	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.006 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	40487	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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: LLP

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	0/3449	0.31	0/4675
1	B	0.15	0/3427	0.32	0/4644
1	C	0.13	0/3423	0.30	0/4639
1	D	0.15	0/3427	0.30	0/4644
1	E	0.13	0/3405	0.28	0/4612
1	F	0.12	0/3403	0.26	0/4611
All	All	0.14	0/20534	0.29	0/27825

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	3397	3329	3329	9	0
1	B	3377	3313	3314	14	0
1	C	3373	3310	3308	16	0
1	D	3377	3312	3314	11	0
1	E	3356	3297	3299	14	0
1	F	3353	3284	3286	16	0
2	A	106	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	120	0	0	1	0
2	C	59	0	0	1	0
2	D	83	0	0	0	0
2	E	35	0	0	0	0
2	F	6	0	0	0	0
All	All	20642	19845	19850	77	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 2.

All (77) 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:E:146:ALA:O	1:E:147:THR:OG1	1.99	0.78
1:D:338:ARG:NH2	1:D:358:PHE:O	2.18	0.76
1:E:134:ARG:NH1	1:E:140:ARG:O	2.22	0.72
1:C:146:ALA:O	1:C:147:THR:OG1	2.09	0.69
1:B:338:ARG:NH2	1:B:358:PHE:O	2.30	0.65
1:D:350:LEU:HD22	1:D:352:TRP:CD1	2.32	0.64
1:F:144:ARG:NH2	1:F:158:SER:O	2.30	0.64
1:B:117:ARG:HG3	1:B:290:LEU:HD13	1.82	0.61
1:B:144:ARG:NH2	1:B:158:SER:O	2.31	0.61
1:C:117:ARG:HG3	1:C:290:LEU:HD13	1.83	0.61
1:B:301:GLN:OE1	1:D:33:THR:HG23	2.00	0.61
1:D:144:ARG:NH2	1:D:158:SER:O	2.35	0.60
1:F:373:ASP:OD1	1:F:375:GLY:N	2.31	0.60
1:A:134:ARG:NH1	1:A:140:ARG:O	2.37	0.57
1:E:338:ARG:NH2	1:E:358:PHE:O	2.38	0.56
1:B:15:VAL:HG12	1:B:39:HIS:CE1	2.42	0.55
1:D:15:VAL:HG12	1:D:39:HIS:CE1	2.42	0.55
1:F:146:ALA:O	1:F:147:THR:OG1	2.16	0.54
1:B:1:GLY:N	2:B:502:HOH:O	2.40	0.54
1:E:333:LEU:HB3	1:E:412:LEU:HD13	1.90	0.54
1:C:140:ARG:NH1	1:C:179:ASN:OD1	2.41	0.53
1:B:175:GLU:O	1:B:179:ASN:ND2	2.42	0.53
1:E:15:VAL:HG12	1:E:39:HIS:CE1	2.44	0.52
1:A:209:PRO:HB2	1:A:254:MET:HG2	1.92	0.51
1:C:15:VAL:HG12	1:C:39:HIS:CE1	2.46	0.51
1:F:112:SER:OG	1:F:255:GLN:OE1	2.14	0.50
1:B:209:PRO:HB2	1:B:254:MET:HG2	1.93	0.50
1:A:136:PHE:O	1:A:296:ARG:NH1	2.46	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:144:ARG:NH2	1:C:158:SER:O	2.44	0.49
1:F:335:GLN:OE1	1:F:338:ARG:NH1	2.44	0.49
1:F:116:LEU:C	1:F:116:LEU:HD12	2.37	0.48
1:B:301:GLN:CD	1:D:33:THR:HG23	2.39	0.48
1:E:154:ASP:N	1:E:154:ASP:OD1	2.46	0.48
1:C:365:PRO:HG2	1:C:389:MET:HE1	1.96	0.48
1:F:188:ILE:HD11	1:F:226:GLU:HG2	1.97	0.47
1:F:115:LYS:NZ	1:F:284:GLU:OE2	2.48	0.47
1:D:147:THR:HA	1:D:160:TYR:O	2.15	0.47
1:D:394:ARG:NH1	1:D:404:ILE:O	2.47	0.47
1:A:190:GLY:O	1:A:194:VAL:HG23	2.15	0.47
1:C:386:ASP:OD1	1:C:387:VAL:N	2.47	0.47
1:D:116:LEU:HD12	1:D:116:LEU:C	2.39	0.46
1:E:116:LEU:HD12	1:E:116:LEU:C	2.39	0.46
1:A:333:LEU:HD22	1:A:421:VAL:HG21	1.98	0.46
1:E:211:GLY:HA2	1:E:219:VAL:HG12	1.96	0.46
1:E:192:ARG:HG2	1:E:230:LEU:HD11	1.98	0.46
1:B:290:LEU:O	1:B:297:LYS:NZ	2.49	0.46
1:B:136:PHE:O	1:B:296:ARG:NH1	2.49	0.45
1:F:126:THR:HG21	1:F:150:HIS:HB3	1.99	0.45
1:A:12:GLU:O	1:A:15:VAL:HG12	2.17	0.45
1:F:211:GLY:HA2	1:F:219:VAL:HG12	1.99	0.45
1:F:333:LEU:CD2	1:F:421:VAL:HG21	2.47	0.45
1:B:11:THR:O	1:B:15:VAL:HG13	2.17	0.44
1:E:63:PHE:CG	1:E:405:GLY:HA3	2.52	0.44
1:D:117:ARG:HG3	1:D:290:LEU:HD13	1.99	0.44
1:A:147:THR:HA	1:A:160:TYR:O	2.19	0.43
1:C:295:ASP:N	1:C:295:ASP:OD1	2.51	0.43
1:B:159:GLY:HA3	1:B:182:LEU:HD11	1.99	0.43
1:C:22:ARG:NH2	1:C:38:PRO:O	2.50	0.43
1:C:134:ARG:NH1	1:C:140:ARG:O	2.45	0.43
1:F:290:LEU:O	1:F:297:LYS:NZ	2.50	0.42
1:C:116:LEU:C	1:C:116:LEU:HD12	2.44	0.42
1:C:190:GLY:O	1:C:194:VAL:HG23	2.20	0.42
1:E:147:THR:HA	1:E:160:TYR:O	2.20	0.42
1:E:333:LEU:HD22	1:E:421:VAL:HG21	2.01	0.42
1:C:34:ARG:NH1	2:C:501:HOH:O	2.47	0.42
1:E:379:ARG:NH1	1:E:381:GLU:OE1	2.52	0.42
1:D:362:HIS:ND1	1:D:409:SER:OG	2.38	0.41
1:C:285:ASP:OD1	1:C:285:ASP:N	2.53	0.41
1:F:188:ILE:HD12	1:F:191:MET:HE2	2.03	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:LEU:HB3	1:A:412:LEU:HD13	2.02	0.41
1:B:98:LEU:HD12	1:B:308:PRO:HB3	2.03	0.41
1:F:394:ARG:NH2	1:F:404:ILE:O	2.54	0.41
1:C:96:HIS:O	1:C:99:GLU:HG2	2.22	0.40
1:E:285:ASP:N	1:E:285:ASP:OD1	2.53	0.40
1:F:404:ILE:HD12	1:F:408:GLY:HA2	2.03	0.40
1:F:136:PHE:O	1:F:296:ARG:NH1	2.54	0.40
1:A:30:THR:HA	1:C:117:ARG:HD3	2.02	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	440/464 (95%)	427 (97%)	12 (3%)	1 (0%)	43	48
1	B	438/464 (94%)	428 (98%)	9 (2%)	1 (0%)	43	48
1	C	437/464 (94%)	424 (97%)	13 (3%)	0	100	100
1	D	438/464 (94%)	422 (96%)	16 (4%)	0	100	100
1	E	433/464 (93%)	420 (97%)	13 (3%)	0	100	100
1	F	435/464 (94%)	419 (96%)	16 (4%)	0	100	100
All	All	2621/2784 (94%)	2540 (97%)	79 (3%)	2 (0%)	48	54

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	406	GLY
1	A	408	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	344/363 (95%)	337 (98%)	7 (2%)	48	58
1	B	342/363 (94%)	331 (97%)	11 (3%)	34	40
1	C	342/363 (94%)	335 (98%)	7 (2%)	48	58
1	D	342/363 (94%)	334 (98%)	8 (2%)	44	53
1	E	340/363 (94%)	333 (98%)	7 (2%)	47	56
1	F	339/363 (93%)	334 (98%)	5 (2%)	57	66
All	All	2049/2178 (94%)	2004 (98%)	45 (2%)	45	54

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	23	SER
1	A	301	GLN
1	A	333	LEU
1	A	373	ASP
1	A	389	MET
1	A	428	PHE
1	B	15	VAL
1	B	19	GLN
1	B	115	LYS
1	B	301	GLN
1	B	325	ASP
1	B	333	LEU
1	B	350	LEU
1	B	372	THR
1	B	373	ASP
1	B	407	ARG
1	B	428	PHE
1	C	2	SER
1	C	15	VAL
1	C	292	ARG
1	C	333	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	350	LEU
1	C	376	SER
1	C	428	PHE
1	D	15	VAL
1	D	188	ILE
1	D	226	GLU
1	D	301	GLN
1	D	333	LEU
1	D	350	LEU
1	D	388	LYS
1	D	428	PHE
1	E	15	VAL
1	E	95	SER
1	E	154	ASP
1	E	325	ASP
1	E	333	LEU
1	E	350	LEU
1	E	373	ASP
1	F	226	GLU
1	F	325	ASP
1	F	333	LEU
1	F	389	MET
1	F	428	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	HIS
1	A	179	ASN
1	A	197	GLN
1	A	261	GLN
1	C	47	GLN
1	C	50	HIS
1	D	57	ASN
1	D	167	GLN
1	D	197	GLN
1	E	47	GLN
1	E	198	HIS
1	F	233	GLN
1	F	343	HIS

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
1	LLP	D	269	1	23,24,25	2.58	6 (26%)	25,32,34	1.27	3 (12%)
1	LLP	B	269	1	23,24,25	2.54	6 (26%)	25,32,34	1.39	3 (12%)
1	LLP	E	269	1	23,24,25	2.58	6 (26%)	25,32,34	1.26	3 (12%)
1	LLP	F	269	1	23,24,25	2.57	6 (26%)	25,32,34	1.16	2 (8%)
1	LLP	A	269	1	23,24,25	2.58	6 (26%)	25,32,34	1.35	3 (12%)
1	LLP	C	269	1	23,24,25	2.57	6 (26%)	25,32,34	1.30	2 (8%)

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
1	LLP	D	269	1	-	5/16/17/19	0/1/1/1
1	LLP	B	269	1	-	6/16/17/19	0/1/1/1
1	LLP	E	269	1	-	10/16/17/19	0/1/1/1
1	LLP	F	269	1	-	5/16/17/19	0/1/1/1
1	LLP	A	269	1	-	6/16/17/19	0/1/1/1
1	LLP	C	269	1	-	4/16/17/19	0/1/1/1

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	269	LLP	C4-C4'	7.77	1.63	1.46

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	269	LLP	C4-C4'	7.73	1.63	1.46
1	C	269	LLP	C4-C4'	7.67	1.63	1.46
1	D	269	LLP	C4-C4'	7.67	1.63	1.46
1	B	269	LLP	C4-C4'	7.61	1.62	1.46
1	F	269	LLP	C4-C4'	7.56	1.62	1.46
1	D	269	LLP	C4'-NZ	5.50	1.45	1.27
1	A	269	LLP	C4'-NZ	5.39	1.45	1.27
1	C	269	LLP	C4'-NZ	5.38	1.45	1.27
1	E	269	LLP	C4'-NZ	5.36	1.45	1.27
1	B	269	LLP	C4'-NZ	5.34	1.45	1.27
1	F	269	LLP	C4'-NZ	5.19	1.44	1.27
1	B	269	LLP	C2'-C2	4.03	1.56	1.50
1	A	269	LLP	C2'-C2	3.92	1.56	1.50
1	C	269	LLP	C2'-C2	3.87	1.56	1.50
1	D	269	LLP	C2'-C2	3.81	1.56	1.50
1	E	269	LLP	C2'-C2	3.81	1.56	1.50
1	F	269	LLP	C2'-C2	3.77	1.56	1.50
1	F	269	LLP	C4-C5	-3.58	1.37	1.42
1	E	269	LLP	C4-C5	-3.34	1.37	1.42
1	A	269	LLP	C4-C5	-3.24	1.37	1.42
1	B	269	LLP	C4-C5	-3.22	1.37	1.42
1	C	269	LLP	C4-C5	-3.22	1.37	1.42
1	F	269	LLP	C6-N1	3.21	1.40	1.34
1	D	269	LLP	C4-C5	-3.18	1.37	1.42
1	E	269	LLP	C6-N1	3.14	1.40	1.34
1	C	269	LLP	C6-N1	3.07	1.40	1.34
1	D	269	LLP	C6-N1	3.06	1.40	1.34
1	A	269	LLP	C6-N1	3.05	1.40	1.34
1	B	269	LLP	C6-N1	2.92	1.40	1.34
1	C	269	LLP	C5'-C5	2.24	1.56	1.50
1	E	269	LLP	C5'-C5	2.18	1.56	1.50
1	F	269	LLP	C5'-C5	2.17	1.56	1.50
1	D	269	LLP	C5'-C5	2.16	1.56	1.50
1	A	269	LLP	C5'-C5	2.14	1.56	1.50
1	B	269	LLP	C5'-C5	2.09	1.56	1.50

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	269	LLP	C4-C4'-NZ	-3.41	108.29	124.04
1	F	269	LLP	C4-C4'-NZ	-3.38	108.44	124.04
1	C	269	LLP	C4-C4'-NZ	-3.29	108.85	124.04

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	LLP	C4-C4'-NZ	-3.29	108.86	124.04
1	D	269	LLP	C4-C4'-NZ	-3.23	109.14	124.04
1	E	269	LLP	C4-C4'-NZ	-3.13	109.60	124.04
1	B	269	LLP	C4-C3-C2	3.04	121.86	120.14
1	A	269	LLP	C4-C3-C2	2.91	121.78	120.14
1	C	269	LLP	C4-C3-C2	2.74	121.69	120.14
1	D	269	LLP	C4-C3-C2	2.46	121.53	120.14
1	A	269	LLP	OP4-C5'-C5	2.41	113.87	109.36
1	E	269	LLP	C4-C3-C2	2.40	121.49	120.14
1	B	269	LLP	OP4-C5'-C5	2.29	113.64	109.36
1	E	269	LLP	OP4-C5'-C5	2.25	113.57	109.36
1	D	269	LLP	OP4-C5'-C5	2.22	113.52	109.36
1	F	269	LLP	C5-C6-N1	-2.16	120.32	123.83

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	269	LLP	O-C-CA-CB
1	A	269	LLP	CG-CD-CE-NZ
1	B	269	LLP	O-C-CA-CB
1	C	269	LLP	O-C-CA-CB
1	D	269	LLP	C4-C4'-NZ-CE
1	D	269	LLP	O-C-CA-CB
1	D	269	LLP	CG-CD-CE-NZ
1	E	269	LLP	C4-C5-C5'-OP4
1	E	269	LLP	C6-C5-C5'-OP4
1	E	269	LLP	C5'-OP4-P-OP2
1	E	269	LLP	C5'-OP4-P-OP3
1	E	269	LLP	O-C-CA-CB
1	F	269	LLP	O-C-CA-CB
1	A	269	LLP	C4-C4'-NZ-CE
1	C	269	LLP	C4-C4'-NZ-CE
1	E	269	LLP	C4-C4'-NZ-CE
1	B	269	LLP	C4-C4'-NZ-CE
1	F	269	LLP	C4-C4'-NZ-CE
1	B	269	LLP	CG-CD-CE-NZ
1	C	269	LLP	CG-CD-CE-NZ
1	E	269	LLP	CG-CD-CE-NZ
1	F	269	LLP	C3-C4-C4'-NZ
1	E	269	LLP	C5'-OP4-P-OP1
1	B	269	LLP	C3-C4-C4'-NZ

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
1	D	269	LLP	C3-C4-C4'-NZ
1	E	269	LLP	CD-CE-NZ-C4'
1	A	269	LLP	C3-C4-C4'-NZ
1	E	269	LLP	C3-C4-C4'-NZ
1	F	269	LLP	CD-CE-NZ-C4'
1	C	269	LLP	C3-C4-C4'-NZ
1	A	269	LLP	C5'-OP4-P-OP2
1	B	269	LLP	C5'-OP4-P-OP2
1	A	269	LLP	CD-CE-NZ-C4'
1	D	269	LLP	CD-CE-NZ-C4'
1	F	269	LLP	C5-C4-C4'-NZ
1	B	269	LLP	CD-CE-NZ-C4'

There are no ring outliers.

No monomer is involved in short contacts.

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

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	442/464 (95%)	-0.19	7 (1%) 70 72	38, 52, 80, 101	0
1	B	440/464 (94%)	-0.12	7 (1%) 70 72	37, 54, 81, 113	0
1	C	439/464 (94%)	0.01	8 (1%) 67 69	36, 60, 94, 142	0
1	D	440/464 (94%)	-0.11	6 (1%) 73 75	34, 57, 85, 123	0
1	E	437/464 (94%)	0.23	12 (2%) 56 57	48, 69, 104, 134	0
1	F	437/464 (94%)	0.59	22 (5%) 34 32	61, 87, 117, 132	0
All	All	2635/2784 (94%)	0.07	62 (2%) 59 61	34, 62, 103, 142	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	171	GLY	5.4
1	C	171	GLY	4.5
1	F	154	ASP	4.3
1	E	164	PHE	4.3
1	E	1	GLY	4.1
1	A	1	GLY	3.9
1	D	441	LEU	3.5
1	C	387	VAL	3.5
1	F	372	THR	3.2
1	E	157	ALA	3.1
1	C	372	THR	2.9
1	F	92	TYR	2.9
1	F	233	GLN	2.9
1	F	33	THR	2.9
1	C	172	VAL	2.8
1	F	1	GLY	2.7
1	F	170	PRO	2.7
1	E	168	PRO	2.7
1	F	157	ALA	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	F	155	PHE	2.7
1	E	158	SER	2.7
1	E	155	PHE	2.6
1	B	189	GLU	2.6
1	B	383	ALA	2.6
1	F	196	ALA	2.6
1	E	441	LEU	2.5
1	D	384	ARG	2.5
1	B	1	GLY	2.5
1	B	384	ARG	2.5
1	D	93	ALA	2.5
1	F	383	ALA	2.5
1	E	154	ASP	2.5
1	B	155	PHE	2.4
1	D	1	GLY	2.4
1	C	407	ARG	2.4
1	F	131	ARG	2.4
1	F	438	GLU	2.4
1	A	92	TYR	2.3
1	B	154	ASP	2.3
1	D	92	TYR	2.3
1	F	200	SER	2.3
1	A	171	GLY	2.3
1	C	384	ARG	2.3
1	C	156	SER	2.3
1	A	37	ASP	2.2
1	F	350	LEU	2.2
1	E	384	ARG	2.2
1	A	293	GLY	2.2
1	F	148	HIS	2.2
1	B	372	THR	2.2
1	A	16	PHE	2.2
1	F	366	GLY	2.2
1	F	171	GLY	2.2
1	F	232	ARG	2.1
1	F	164	PHE	2.1
1	E	170	PRO	2.1
1	F	370	ASN	2.1
1	A	443	TRP	2.1
1	D	159	GLY	2.1
1	F	387	VAL	2.1
1	C	155	PHE	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	199	GLY	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
1	LLP	F	269	24/25	0.89	0.16	65,82,100,101	0
1	LLP	E	269	24/25	0.92	0.12	54,69,80,83	0
1	LLP	C	269	24/25	0.93	0.12	46,59,72,72	0
1	LLP	D	269	24/25	0.94	0.10	40,54,63,64	0
1	LLP	A	269	24/25	0.95	0.09	36,52,62,63	0
1	LLP	B	269	24/25	0.95	0.08	37,50,60,62	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.