



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 5, 2026 – 01:10 PM UTC

PDB ID : 3UNV / pdb_00003unv
Title : Pantoea agglomerans Phenylalanine Aminomutase
Authors : Geiger, J.; Strom, S.
Deposited on : 2011-11-16
Resolution : 1.54 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

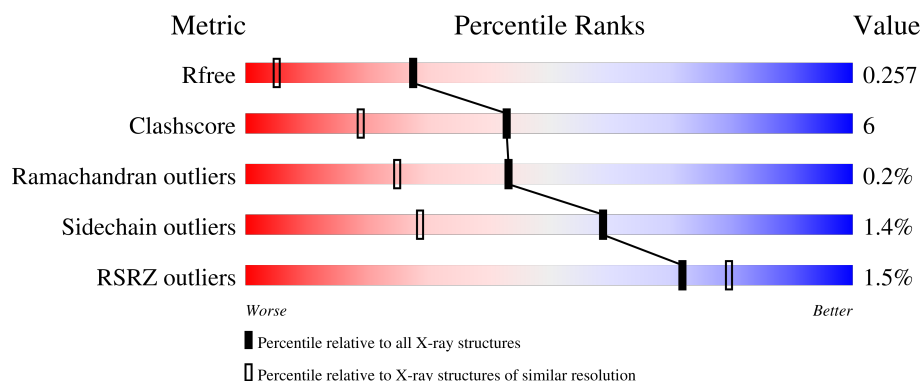
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.54 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	1003 (1.54-1.54)
Clashscore	190562	1025 (1.54-1.54)
Ramachandran outliers	187476	1007 (1.54-1.54)
Sidechain outliers	187428	1007 (1.54-1.54)
RSRZ outliers	180081	1002 (1.54-1.54)

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	547	 76% 16% • 6%
1	B	547	 79% 14% • 6%

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
5	EDO	B	1003	-	-	X	-
5	EDO	B	1004	-	-	X	-
6	PO4	A	550	-	X	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8538 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.

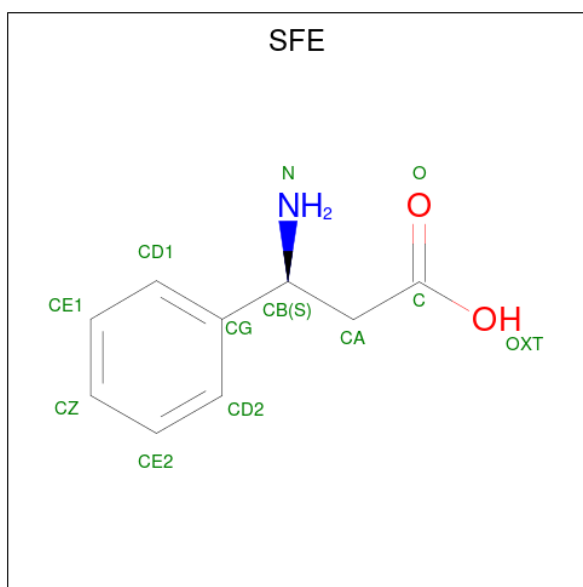
- Molecule 1 is a protein called AdmH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	514	Total	C	N	O	S	1	3	0
			3981	2494	692	770	25			
1	B	514	Total	C	N	O	S	7	3	0
			3970	2489	689	767	25			

There are 16 discrepancies between the modelled and reference sequences:

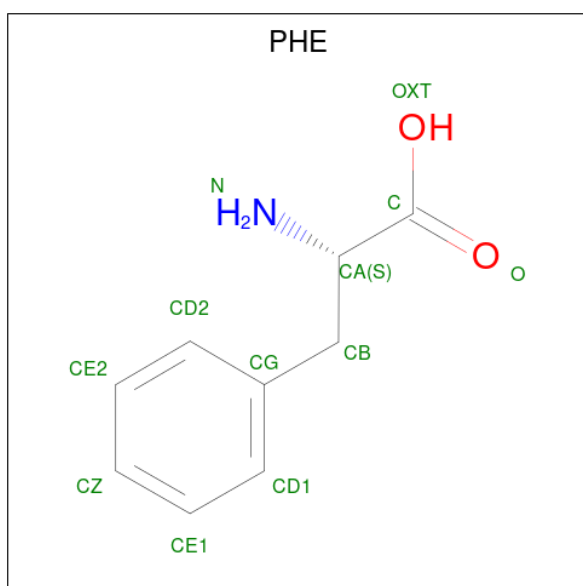
Chain	Residue	Modelled	Actual	Comment	Reference
A	542	LEU	-	expression tag	UNP Q84FL5
A	543	GLU	-	expression tag	UNP Q84FL5
A	544	HIS	-	expression tag	UNP Q84FL5
A	545	HIS	-	expression tag	UNP Q84FL5
A	546	HIS	-	expression tag	UNP Q84FL5
A	547	HIS	-	expression tag	UNP Q84FL5
A	548	HIS	-	expression tag	UNP Q84FL5
A	549	HIS	-	expression tag	UNP Q84FL5
B	542	LEU	-	expression tag	UNP Q84FL5
B	543	GLU	-	expression tag	UNP Q84FL5
B	544	HIS	-	expression tag	UNP Q84FL5
B	545	HIS	-	expression tag	UNP Q84FL5
B	546	HIS	-	expression tag	UNP Q84FL5
B	547	HIS	-	expression tag	UNP Q84FL5
B	548	HIS	-	expression tag	UNP Q84FL5
B	549	HIS	-	expression tag	UNP Q84FL5

- Molecule 2 is (3S)-3-amino-3-phenylpropanoic acid (CCD ID: SFE) (formula: C₉H₁₁NO₂).



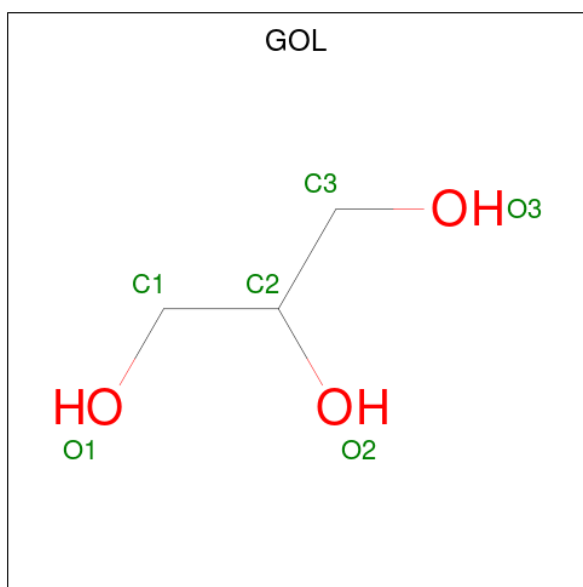
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	1
			12	9	1	2		
2	B	1	Total	C	N	O	0	0
			12	9	1	2		

- Molecule 3 is PHENYLALANINE (CCD ID: PHE) (formula: C₉H₁₁NO₂).



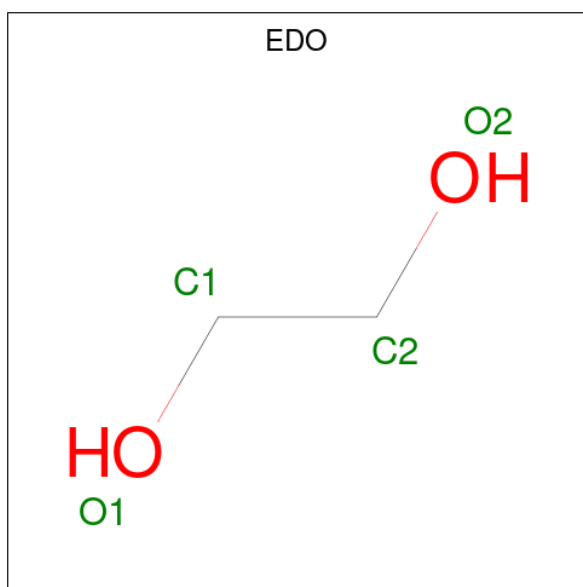
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	1
			12	9	1	2		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



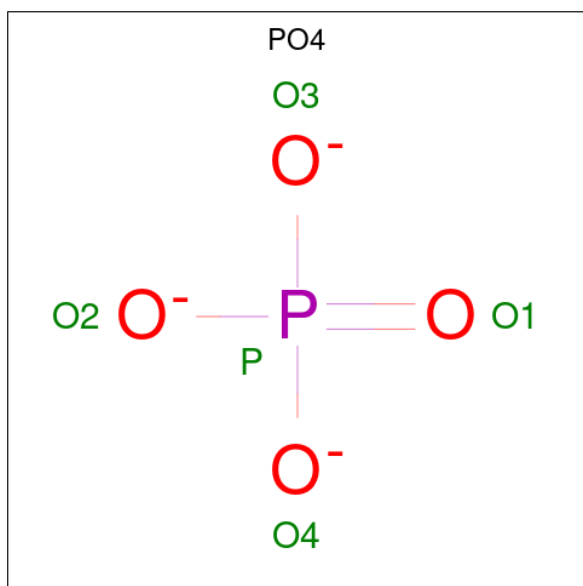
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 6 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	P	0	0
			5	4	1		
6	A	1	Total	O	P	0	0
			5	4	1		

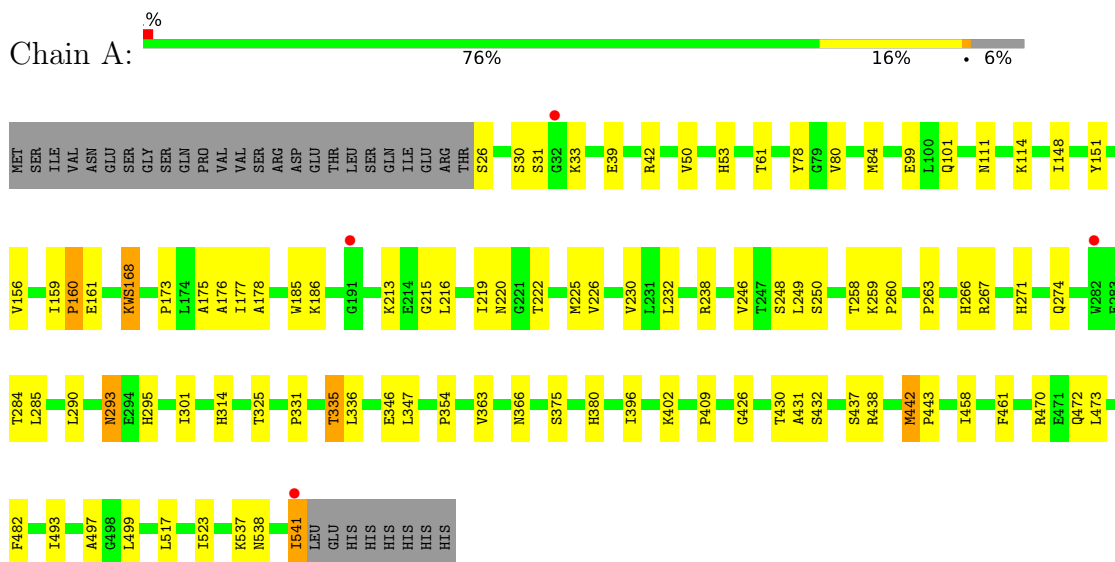
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	243	Total 243	O 243	0	0
7	B	248	Total 248	O 248	0	0

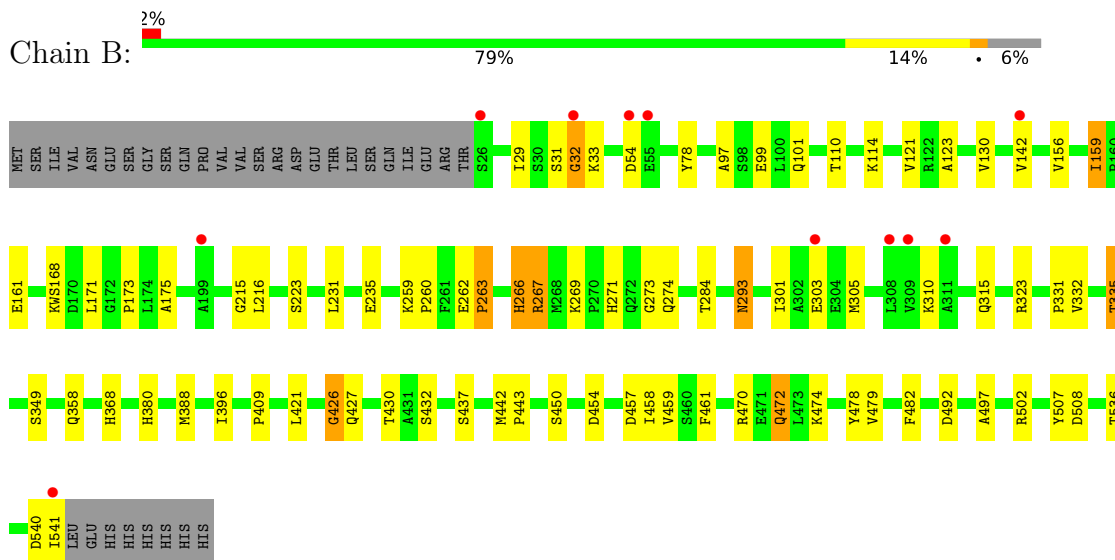
3 Residue-property plots [i](#)

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: AdmH



• Molecule 1: AdmH



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	153.96Å 185.84Å 72.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.54 50.00 – 1.54	Depositor EDS
% Data completeness (in resolution range)	91.1 (50.00-1.54) 91.2 (50.00-1.54)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 1.54Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.204 , 0.247 0.212 , 0.257	Depositor DCC
R_{free} test set	7034 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	16.0	Xtriage
Anisotropy	0.658	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8538	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% 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: EDO, PO4, KWS, GOL, SFE

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	1.57	23/4016 (0.6%)	1.34	20/5433 (0.4%)
1	B	1.59	28/4023 (0.7%)	1.30	14/5444 (0.3%)
All	All	1.58	51/8039 (0.6%)	1.32	34/10877 (0.3%)

The worst 5 of 51 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	409	PRO	CA-C	9.38	1.57	1.51
1	A	336	LEU	N-CA	9.06	1.57	1.46
1	B	332	VAL	CA-CB	-7.59	1.45	1.54
1	B	156	VAL	CA-CB	7.09	1.60	1.54
1	B	421	LEU	N-CA	6.73	1.54	1.46

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	346	GLU	N-CA-C	-7.88	102.00	111.69
1	B	426	GLY	N-CA-C	-7.17	103.77	113.37
1	A	80	VAL	N-CA-C	-6.93	104.97	111.48
1	A	249	LEU	N-CA-C	6.91	118.46	111.07
1	B	223	SER	N-CA-C	6.70	119.42	111.71

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	3981	0	3979	55	0
1	B	3970	0	3978	42	0
2	A	12	0	8	5	0
2	B	12	0	9	1	0
3	A	12	0	8	1	0
4	A	6	0	8	1	0
4	B	12	0	16	1	0
5	A	16	0	24	5	0
5	B	16	0	24	11	0
6	A	10	0	0	1	0
7	A	243	0	0	6	0
7	B	248	0	0	2	0
All	All	8538	0	8054	99	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.

The worst 5 of 99 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)
7:A:779:HOH:O	5:B:550:EDO:H22	1.59	1.01
1:B:454:ASP:OD1	5:B:1003:EDO:H21	1.68	0.92
1:A:538:ASN:O	1:A:541:ILE:HG22	1.81	0.81
1:B:266:HIS:HD2	1:B:274:GLN:HE21	1.33	0.77
1:A:266:HIS:HD2	1:A:274:GLN:HE21	1.32	0.74

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	511/547 (93%)	502 (98%)	9 (2%)	0	100	100
1	B	512/547 (94%)	500 (98%)	10 (2%)	2 (0%)	30	12
All	All	1023/1094 (94%)	1002 (98%)	19 (2%)	2 (0%)	43	25

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	33	LYS
1	B	32	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	437/469 (93%)	429 (98%)	8 (2%)	51	21
1	B	436/469 (93%)	432 (99%)	4 (1%)	70	46
All	All	873/938 (93%)	861 (99%)	12 (1%)	59	29

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	541	ILE
1	B	263	PRO
1	B	540	ASP
1	B	293	ASN
1	A	293	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 30 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	472	GLN
1	B	393	ASN
1	B	102	ASN
1	B	472	GLN

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Mol	Chain	Res	Type
1	B	295	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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	KWS	A	168[A]	1,2	13,15,16	3.56	4 (30%)	17,21,23	5.70	10 (58%)
1	KWS	A	168[B]	3,1	13,15,16	3.95	4 (30%)	17,21,23	4.91	10 (58%)
1	KWS	B	168	1,2	13,15,16	3.68	3 (23%)	17,21,23	4.71	8 (47%)

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	KWS	A	168[A]	1,2	-	2/8/27/28	0/1/1/1
1	KWS	A	168[B]	3,1	-	3/8/27/28	0/1/1/1
1	KWS	B	168	1,2	-	3/8/27/28	0/1/1/1

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	168	KWS	O2-C2	11.98	1.47	1.23
1	A	168[B]	KWS	O2-C2	11.68	1.46	1.23
1	A	168[A]	KWS	O2-C2	11.21	1.45	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	168[B]	KWS	CA1-N1	5.45	1.62	1.47
1	A	168[B]	KWS	C2-N3	-4.09	1.30	1.40

The worst 5 of 28 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	168[A]	KWS	CA2-C2-N3	14.33	115.53	103.50
1	B	168	KWS	CA2-C2-N3	12.23	113.77	103.50
1	A	168[A]	KWS	C2-N3-C1	-11.62	102.69	108.07
1	A	168[A]	KWS	C2-CA2-N2	-11.01	101.06	108.95
1	A	168[B]	KWS	CA2-C2-N3	10.91	112.66	103.50

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	168[A]	KWS	N2-C1-CA1-CB
1	A	168[A]	KWS	N3-C1-CA1-CB
1	A	168[B]	KWS	N2-C1-CA1-CB
1	A	168[B]	KWS	N3-C1-CA1-CB
1	B	168	KWS	N2-C1-CA1-CB

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	168[A]	KWS	3	0
1	A	168[B]	KWS	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

16 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	GOL	A	1002	-	5,5,5	0.68	0	5,5,5	0.35	0
4	GOL	B	1002	-	5,5,5	0.23	0	5,5,5	0.90	0
5	EDO	A	1004	-	3,3,3	0.69	0	2,2,2	0.84	0
5	EDO	A	1005	-	3,3,3	0.38	0	2,2,2	0.84	0
6	PO4	A	550	-	4,4,4	2.45	3 (75%)	6,6,6	1.58	1 (16%)
4	GOL	B	1001	-	5,5,5	0.98	0	5,5,5	0.69	0
6	PO4	A	1008	-	4,4,4	0.92	0	6,6,6	1.07	1 (16%)
2	SFE	A	1000[A]	1	11,12,12	0.96	0	13,15,15	1.62	2 (15%)
2	SFE	B	1000	1	11,12,12	1.05	1 (9%)	13,15,15	1.45	3 (23%)
5	EDO	B	1005	-	3,3,3	0.43	0	2,2,2	1.15	0
5	EDO	A	1006	-	3,3,3	0.14	0	2,2,2	1.21	0
5	EDO	A	1007	-	3,3,3	0.56	0	2,2,2	0.34	0
5	EDO	B	1004	-	3,3,3	0.39	0	2,2,2	1.16	0
5	EDO	B	1003	-	3,3,3	1.00	0	2,2,2	0.73	0
3	PHE	A	1001[B]	1	11,12,12	0.78	0	11,15,15	1.17	1 (9%)
5	EDO	B	550	-	3,3,3	0.17	0	2,2,2	0.81	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	GOL	A	1002	-	-	0/4/4/4	-
4	GOL	B	1002	-	-	0/4/4/4	-
5	EDO	A	1004	-	-	0/1/1/1	-
5	EDO	A	1005	-	-	1/1/1/1	-
4	GOL	B	1001	-	-	0/4/4/4	-
2	SFE	A	1000[A]	1	-	2/8/8/8	0/1/1/1
2	SFE	B	1000	1	-	1/8/8/8	0/1/1/1
5	EDO	B	1005	-	-	1/1/1/1	-
5	EDO	A	1006	-	-	0/1/1/1	-
5	EDO	A	1007	-	-	0/1/1/1	-
5	EDO	B	1004	-	-	1/1/1/1	-
5	EDO	B	1003	-	-	0/1/1/1	-
3	PHE	A	1001[B]	1	-	6/8/8/8	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	B	550	-	-	1/1/1/1	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	550	PO4	P-O3	-3.45	1.44	1.54
6	A	550	PO4	P-O1	-2.51	1.44	1.50
6	A	550	PO4	P-O4	-2.33	1.47	1.54
2	B	1000	SFE	OXT-C	-2.09	1.23	1.30

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1000[A]	SFE	CA-CB-CG	-3.37	104.94	111.34
2	A	1000[A]	SFE	CD1-CG-CD2	2.95	121.96	118.30
2	B	1000	SFE	CE2-CD2-CG	-2.90	117.18	120.65
6	A	550	PO4	O3-P-O1	2.89	121.18	110.95
3	A	1001[B]	PHE	CG-CB-CA	-2.85	108.31	114.13

There are no chirality outliers.

5 of 13 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1001[B]	PHE	O-C-CA-N
3	A	1001[B]	PHE	OXT-C-CA-N
3	A	1001[B]	PHE	CA-CB-CG-CD1
3	A	1001[B]	PHE	CA-CB-CG-CD2
5	A	1005	EDO	O1-C1-C2-O2

There are no ring outliers.

12 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1002	GOL	1	0
5	A	1005	EDO	3	0
4	B	1001	GOL	1	0
6	A	1008	PO4	1	0
2	A	1000[A]	SFE	5	0
2	B	1000	SFE	1	0
5	A	1006	EDO	1	0
5	A	1007	EDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1004	EDO	4	0
5	B	1003	EDO	5	0
3	A	1001[B]	PHE	1	0
5	B	550	EDO	2	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 [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	513/547 (93%)	-0.06	4 (0%) 82 87	9, 19, 32, 42	3 (0%)
1	B	513/547 (93%)	0.08	11 (2%) 63 73	8, 20, 33, 44	5 (0%)
All	All	1026/1094 (93%)	0.01	15 (1%) 72 79	8, 20, 33, 44	8 (0%)

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	541	ILE	8.1
1	B	541	ILE	8.0
1	B	54	ASP	4.8
1	B	32	GLY	3.4
1	B	55	GLU	3.0

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	KWS	A	168[A]	15/16	0.95	0.07	7,13,19,22	15
1	KWS	A	168[B]	15/16	0.95	0.07	19,20,23,27	15
1	KWS	B	168	15/16	0.95	0.07	16,20,24,31	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

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
5	EDO	B	1005	4/4	0.73	0.19	58,59,60,60	0
5	EDO	A	1007	4/4	0.74	0.21	48,53,54,56	0
5	EDO	B	1004	4/4	0.83	0.13	31,37,39,45	0
5	EDO	B	1003	4/4	0.83	0.17	29,31,32,32	0
5	EDO	A	1005	4/4	0.84	0.14	41,42,45,45	0
6	PO4	A	1008	5/5	0.86	0.10	53,54,56,56	0
3	PHE	A	1001[B]	12/12	0.87	0.12	24,35,38,38	12
5	EDO	A	1006	4/4	0.89	0.12	29,30,31,33	0
5	EDO	A	1004	4/4	0.89	0.11	24,25,27,28	0
2	SFE	A	1000[A]	12/12	0.89	0.10	23,26,27,27	12
2	SFE	B	1000	12/12	0.91	0.11	24,32,37,38	0
4	GOL	B	1001	6/6	0.93	0.09	22,27,29,29	0
4	GOL	B	1002	6/6	0.93	0.09	26,32,33,37	0
5	EDO	B	550	4/4	0.94	0.08	25,28,30,32	0
6	PO4	A	550	5/5	0.94	0.19	8,10,15,20	5
4	GOL	A	1002	6/6	0.95	0.08	18,24,27,28	0

6.5 Other polymers

There are no such residues in this entry.