



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 05:10 PM UTC

PDB ID : 1XE3 / pdb_00001xe3
Title : Crystal Structure of purine nucleoside phosphorylase DeoD from *Bacillus anthracis*
Authors : Grenha, R.; Levдикov, V.M.; Fogg, M.; Blagova, E.V.; Brannigan, J.A.; Wilkinson, A.J.; Wilson, K.S.; Structural Proteomics in Europe (SPINE)
Deposited on : 2004-09-09
Resolution : 2.24 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
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

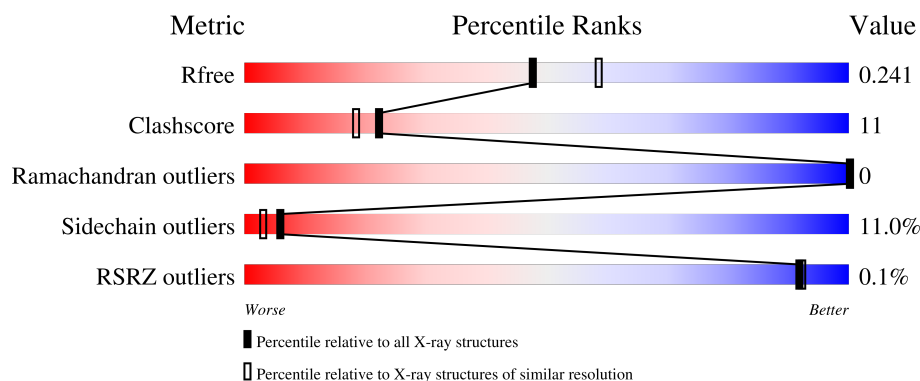
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.24 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	245	 70% 20% . . 5%
1	B	245	 74% 16% 5% 5%
1	C	245	 76% 15% . .
1	D	245	 69% 18% 7% . .
1	E	245	 73% 18% . .

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Mol	Chain	Length	Quality of chain
1	F	245	 A horizontal bar chart showing the quality of chain 1. The bar is divided into three segments: a green segment representing 74%, a yellow segment representing 16%, and a small grey segment representing 5%.

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
2	CL	C	806	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11450 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 purine nucleoside phosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1782	1127	296	347	12			
1	B	233	Total	C	N	O	S	0	0	0
			1780	1127	294	347	12			
1	C	234	Total	C	N	O	S	0	0	0
			1790	1133	297	348	12			
1	D	235	Total	C	N	O	S	0	0	0
			1799	1138	299	350	12			
1	E	234	Total	C	N	O	S	0	0	0
			1789	1132	296	349	12			
1	F	233	Total	C	N	O	S	0	1	0
			1788	1132	297	347	12			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	MET	-	cloning artifact	UNP Q81T09
A	-8	GLY	-	cloning artifact	UNP Q81T09
A	-7	SER	-	cloning artifact	UNP Q81T09
A	-6	SER	-	cloning artifact	UNP Q81T09
A	-5	HIS	-	expression tag	UNP Q81T09
A	-4	HIS	-	expression tag	UNP Q81T09
A	-3	HIS	-	expression tag	UNP Q81T09
A	-2	HIS	-	expression tag	UNP Q81T09
A	-1	HIS	-	expression tag	UNP Q81T09
A	0	HIS	-	expression tag	UNP Q81T09
B	-9	MET	-	cloning artifact	UNP Q81T09
B	-8	GLY	-	cloning artifact	UNP Q81T09
B	-7	SER	-	cloning artifact	UNP Q81T09
B	-6	SER	-	cloning artifact	UNP Q81T09
B	-5	HIS	-	expression tag	UNP Q81T09
B	-4	HIS	-	expression tag	UNP Q81T09
B	-3	HIS	-	expression tag	UNP Q81T09

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	HIS	-	expression tag	UNP Q81T09
B	-1	HIS	-	expression tag	UNP Q81T09
B	0	HIS	-	expression tag	UNP Q81T09
C	-9	MET	-	cloning artifact	UNP Q81T09
C	-8	GLY	-	cloning artifact	UNP Q81T09
C	-7	SER	-	cloning artifact	UNP Q81T09
C	-6	SER	-	cloning artifact	UNP Q81T09
C	-5	HIS	-	expression tag	UNP Q81T09
C	-4	HIS	-	expression tag	UNP Q81T09
C	-3	HIS	-	expression tag	UNP Q81T09
C	-2	HIS	-	expression tag	UNP Q81T09
C	-1	HIS	-	expression tag	UNP Q81T09
C	0	HIS	-	expression tag	UNP Q81T09
D	-9	MET	-	cloning artifact	UNP Q81T09
D	-8	GLY	-	cloning artifact	UNP Q81T09
D	-7	SER	-	cloning artifact	UNP Q81T09
D	-6	SER	-	cloning artifact	UNP Q81T09
D	-5	HIS	-	expression tag	UNP Q81T09
D	-4	HIS	-	expression tag	UNP Q81T09
D	-3	HIS	-	expression tag	UNP Q81T09
D	-2	HIS	-	expression tag	UNP Q81T09
D	-1	HIS	-	expression tag	UNP Q81T09
D	0	HIS	-	expression tag	UNP Q81T09
E	-9	MET	-	cloning artifact	UNP Q81T09
E	-8	GLY	-	cloning artifact	UNP Q81T09
E	-7	SER	-	cloning artifact	UNP Q81T09
E	-6	SER	-	cloning artifact	UNP Q81T09
E	-5	HIS	-	expression tag	UNP Q81T09
E	-4	HIS	-	expression tag	UNP Q81T09
E	-3	HIS	-	expression tag	UNP Q81T09
E	-2	HIS	-	expression tag	UNP Q81T09
E	-1	HIS	-	expression tag	UNP Q81T09
E	0	HIS	-	expression tag	UNP Q81T09
F	-9	MET	-	cloning artifact	UNP Q81T09
F	-8	GLY	-	cloning artifact	UNP Q81T09
F	-7	SER	-	cloning artifact	UNP Q81T09
F	-6	SER	-	cloning artifact	UNP Q81T09
F	-5	HIS	-	expression tag	UNP Q81T09
F	-4	HIS	-	expression tag	UNP Q81T09
F	-3	HIS	-	expression tag	UNP Q81T09
F	-2	HIS	-	expression tag	UNP Q81T09
F	-1	HIS	-	expression tag	UNP Q81T09

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	HIS	-	expression tag	UNP Q81T09

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0
2	B	1	Total Cl 1 1	0	0
2	C	2	Total Cl 2 2	0	0
2	D	1	Total Cl 1 1	0	0
2	F	1	Total Cl 1 1	0	0

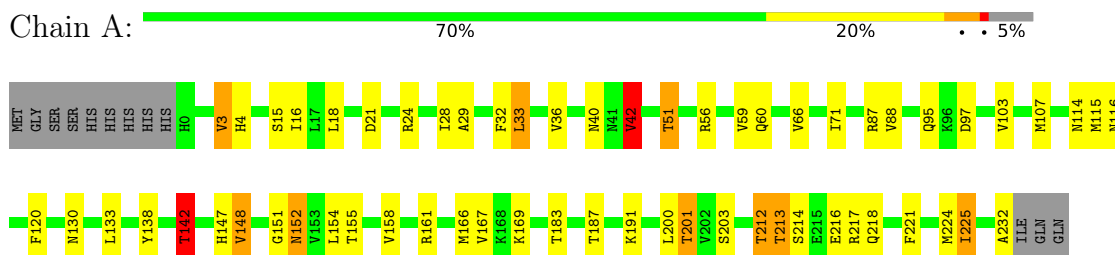
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	148	Total O 148 148	0	0
3	B	114	Total O 114 114	0	0
3	C	115	Total O 115 115	0	0
3	D	115	Total O 115 115	0	0
3	E	122	Total O 122 122	0	0
3	F	102	Total O 102 102	0	0

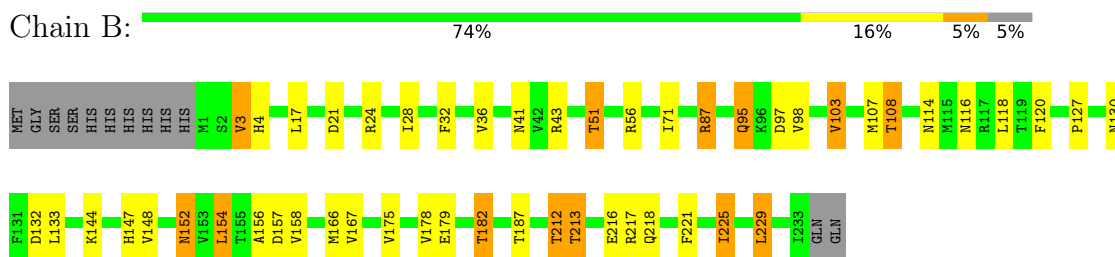
3 Residue-property plots

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 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.

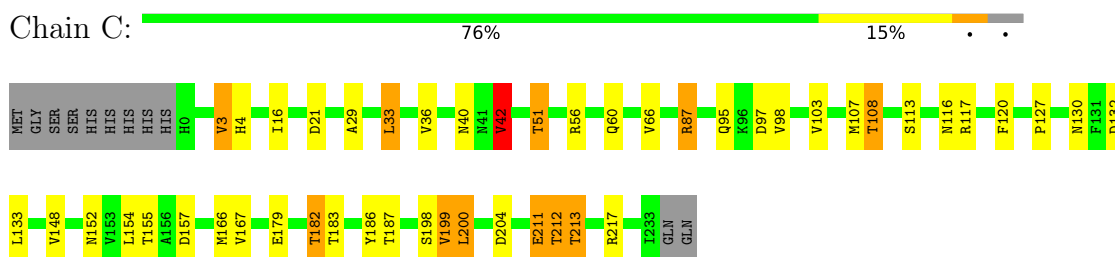
- Molecule 1: purine nucleoside phosphorylase



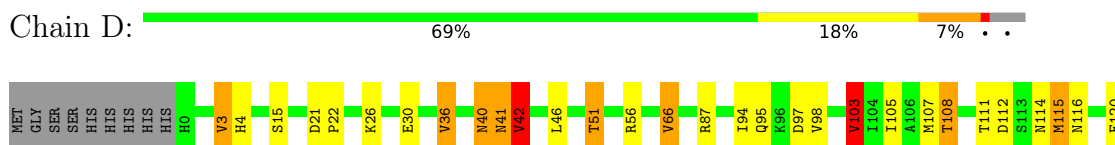
- Molecule 1: purine nucleoside phosphorylase

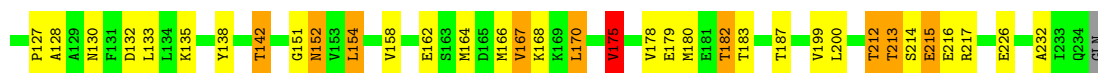


- Molecule 1: purine nucleoside phosphorylase



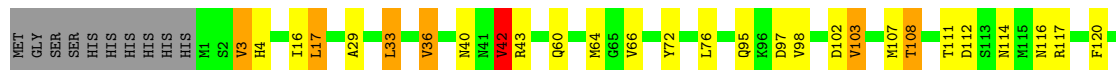
- Molecule 1: purine nucleoside phosphorylase





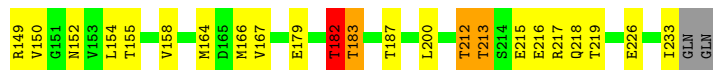
- Molecule 1: purine nucleoside phosphorylase

Chain E: 73% 18%



- Molecule 1: purine nucleoside phosphorylase

Chain F: 74% 16% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.86Å 128.26Å 223.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.24 20.00 – 2.24	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.24) 90.8 (20.00-2.24)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.23Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.181 , 0.235 0.188 , 0.241	Depositor DCC
R_{free} test set	4066 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	22.5	Xtriage
Anisotropy	0.334	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11450	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.90% 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:
CL

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	0.84	4/1809 (0.2%)	0.99	3/2447 (0.1%)
1	B	0.83	1/1806 (0.1%)	0.94	0/2443
1	C	0.85	2/1817 (0.1%)	0.93	2/2458 (0.1%)
1	D	0.90	5/1826 (0.3%)	1.05	2/2470 (0.1%)
1	E	0.88	3/1815 (0.2%)	0.97	4/2455 (0.2%)
1	F	0.87	3/1817 (0.2%)	0.92	0/2457
All	All	0.86	18/10890 (0.2%)	0.97	11/14730 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	66	VAL	CA-CB	11.02	1.59	1.54
1	F	66	VAL	CA-CB	8.73	1.58	1.54
1	C	66	VAL	CA-CB	7.51	1.57	1.54
1	A	66	VAL	CA-CB	7.11	1.57	1.54
1	C	42	VAL	CA-CB	6.78	1.61	1.54

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	111	THR	N-CA-C	6.30	118.89	108.99
1	D	111	THR	N-CA-C	6.30	118.65	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	111	THR	CA-C-N	6.25	129.81	120.31
1	E	111	THR	C-N-CA	6.25	129.81	120.31
1	A	42	VAL	N-CA-CB	5.65	117.70	110.13

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	40	ASN	Peptide

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	1782	0	1787	37	0
1	B	1780	0	1791	43	0
1	C	1790	0	1798	41	0
1	D	1799	0	1806	54	0
1	E	1789	0	1799	39	0
1	F	1788	0	1804	40	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	2	0
2	D	1	0	0	0	0
2	F	1	0	0	0	0
3	A	148	0	0	4	0
3	B	114	0	0	5	0
3	C	115	0	0	3	0
3	D	115	0	0	6	0
3	E	122	0	0	2	0
3	F	102	0	0	3	0
All	All	11450	0	10785	228	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 11.

The worst 5 of 228 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:B:51:THR:HG22	1:B:56:ARG:HH11	1.41	0.85
1:A:158:VAL:HB	3:A:931:HOH:O	1.76	0.84
1:A:138:TYR:O	1:A:142:THR:HG23	1.77	0.84
1:E:138:TYR:O	1:E:142:THR:HG22	1.80	0.81
1:B:108:THR:HG22	1:B:127:PRO:O	1.85	0.75

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	231/245 (94%)	231 (100%)	0	0	100	100
1	B	231/245 (94%)	228 (99%)	3 (1%)	0	100	100
1	C	232/245 (95%)	229 (99%)	3 (1%)	0	100	100
1	D	233/245 (95%)	231 (99%)	2 (1%)	0	100	100
1	E	232/245 (95%)	227 (98%)	5 (2%)	0	100	100
1	F	232/245 (95%)	230 (99%)	2 (1%)	0	100	100
All	All	1391/1470 (95%)	1376 (99%)	15 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	192/203 (95%)	170 (88%)	22 (12%)	5	2
1	B	192/203 (95%)	169 (88%)	23 (12%)	5	2
1	C	193/203 (95%)	174 (90%)	19 (10%)	7	4
1	D	194/203 (96%)	172 (89%)	22 (11%)	5	3
1	E	193/203 (95%)	176 (91%)	17 (9%)	9	6
1	F	193/203 (95%)	168 (87%)	25 (13%)	4	2
All	All	1157/1218 (95%)	1029 (89%)	128 (11%)	6	3

5 of 128 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	149	ARG
1	F	182	THR
1	C	42	VAL
1	C	36	VAL
1	F	200	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 37 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	152	ASN
1	F	152	ASN
1	E	218	GLN
1	F	116	ASN
1	C	116	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

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](#)

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(Å²)	Q<0.9	
1	A	233/245 (95%)	-0.41	0	100	100	15, 21, 29, 35	0
1	B	233/245 (95%)	-0.40	0	100	100	16, 21, 29, 35	0
1	C	234/245 (95%)	-0.42	0	100	100	14, 21, 28, 35	0
1	D	235/245 (95%)	-0.37	0	100	100	13, 21, 31, 40	0
1	E	234/245 (95%)	-0.31	1 (0%)	88	90	15, 22, 32, 42	0
1	F	233/245 (95%)	-0.33	0	100	100	11, 22, 33, 40	1 (0%)
All	All	1402/1470 (95%)	-0.37	1 (0%)	92	92	11, 21, 31, 42	1 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	162	GLU	2.3

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($\text{\AA}^2$)	Q<0.9
2	CL	A	805	1/1	0.98	0.13	26,26,26,26	0
2	CL	B	801	1/1	0.98	0.12	23,23,23,23	0
2	CL	C	806	1/1	0.98	0.24	24,24,24,24	0
2	CL	C	802	1/1	0.99	0.05	17,17,17,17	0
2	CL	F	803	1/1	0.99	0.03	17,17,17,17	0
2	CL	D	804	1/1	1.00	0.04	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.