



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:55 PM UTC

PDB ID : 5WZE / pdb_00005wze
Title : The structure of Pseudomonas aeruginosa aminopeptidase PepP
Authors : Bao, R.; Peng, C.T.; Liu, L.; He, L.H.; Li, C.C.; Li, T.; Shen, Y.L.; Zhu, Y.B.;
Song, Y.J.
Deposited on : 2017-01-17
Resolution : 1.78 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

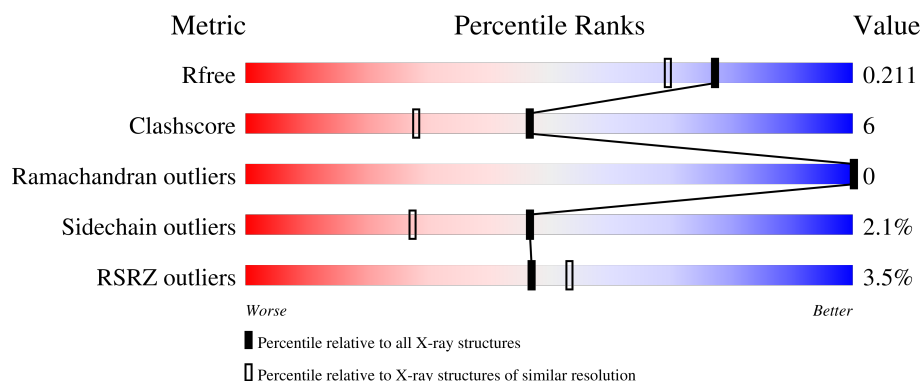
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.78 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	1365 (1.78-1.78)
Clashscore	190562	1395 (1.78-1.78)
Ramachandran outliers	187476	1382 (1.78-1.78)
Sidechain outliers	187428	1382 (1.78-1.78)
RSRZ outliers	180081	1365 (1.78-1.78)

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	452	3% 88% 9% ..
1	B	452	3% 87% 12% .
1	C	452	5% 83% 12% ..
1	D	452	3% 86% 11% ..

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
10	EDO	B	505	-	-	X	-

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 15378 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 Aminopeptidase P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	0	4	0
			3511	2217	627	645	22			
1	B	449	Total	C	N	O	S	0	4	0
			3550	2244	631	654	21			
1	C	442	Total	C	N	O	S	0	7	0
			3523	2229	630	643	21			
1	D	444	Total	C	N	O	S	0	3	0
			3509	2217	624	647	21			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	445	LEU	-	expression tag	UNP Q9HTW6
A	446	GLU	-	expression tag	UNP Q9HTW6
A	447	HIS	-	expression tag	UNP Q9HTW6
A	448	HIS	-	expression tag	UNP Q9HTW6
A	449	HIS	-	expression tag	UNP Q9HTW6
A	450	HIS	-	expression tag	UNP Q9HTW6
A	451	HIS	-	expression tag	UNP Q9HTW6
A	452	HIS	-	expression tag	UNP Q9HTW6
B	445	LEU	-	expression tag	UNP Q9HTW6
B	446	GLU	-	expression tag	UNP Q9HTW6
B	447	HIS	-	expression tag	UNP Q9HTW6
B	448	HIS	-	expression tag	UNP Q9HTW6
B	449	HIS	-	expression tag	UNP Q9HTW6
B	450	HIS	-	expression tag	UNP Q9HTW6
B	451	HIS	-	expression tag	UNP Q9HTW6
B	452	HIS	-	expression tag	UNP Q9HTW6
C	445	LEU	-	expression tag	UNP Q9HTW6
C	446	GLU	-	expression tag	UNP Q9HTW6
C	447	HIS	-	expression tag	UNP Q9HTW6
C	448	HIS	-	expression tag	UNP Q9HTW6
C	449	HIS	-	expression tag	UNP Q9HTW6

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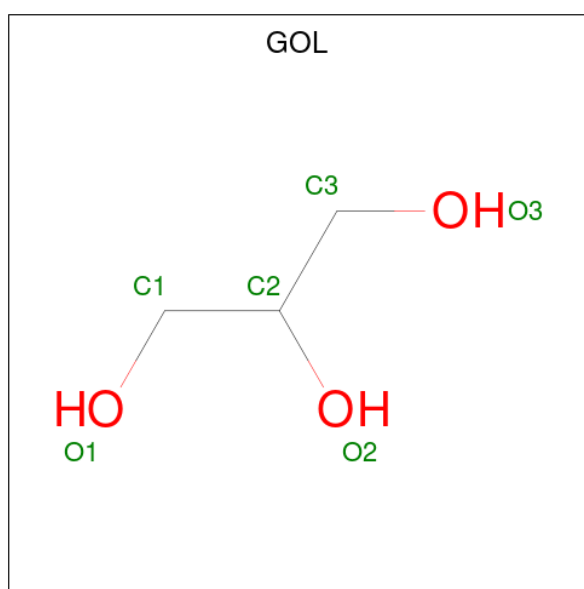
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Chain	Residue	Modelled	Actual	Comment	Reference
C	450	HIS	-	expression tag	UNP Q9HTW6
C	451	HIS	-	expression tag	UNP Q9HTW6
C	452	HIS	-	expression tag	UNP Q9HTW6
D	445	LEU	-	expression tag	UNP Q9HTW6
D	446	GLU	-	expression tag	UNP Q9HTW6
D	447	HIS	-	expression tag	UNP Q9HTW6
D	448	HIS	-	expression tag	UNP Q9HTW6
D	449	HIS	-	expression tag	UNP Q9HTW6
D	450	HIS	-	expression tag	UNP Q9HTW6
D	451	HIS	-	expression tag	UNP Q9HTW6
D	452	HIS	-	expression tag	UNP Q9HTW6

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

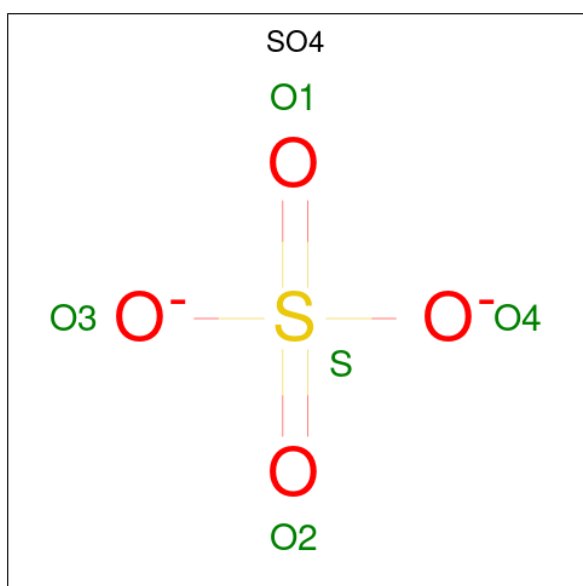
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Na 1 1	0	0
2	B	2	Total Na 2 2	0	0
2	C	1	Total Na 1 1	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



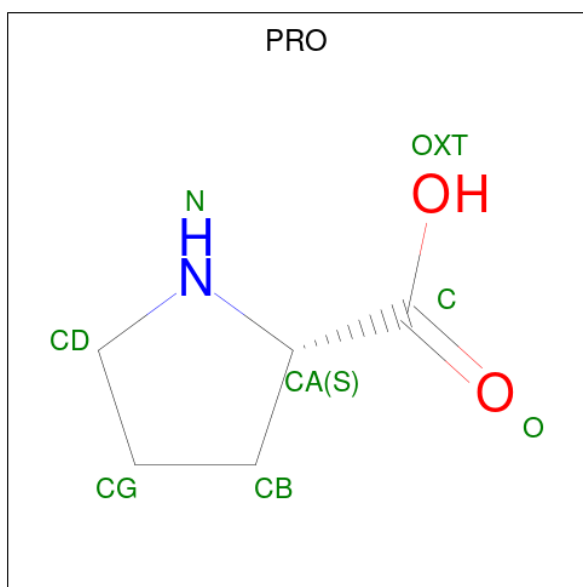
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



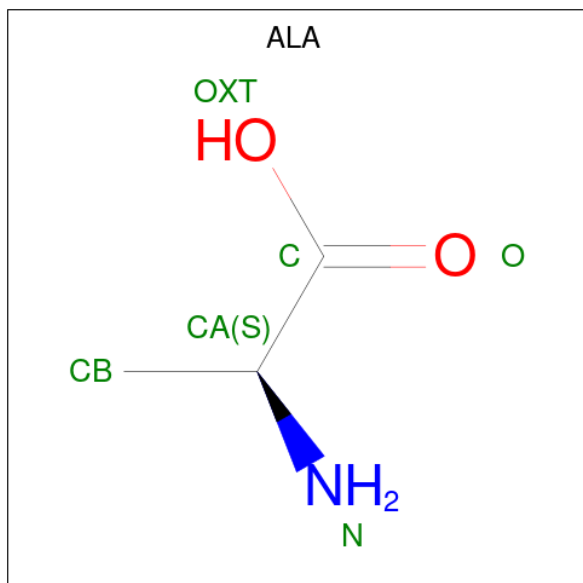
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is PROLINE (CCD ID: PRO) (formula: C₅H₉NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			8	5	1	2		
5	B	1	Total	C	N	O	0	0
			8	5	1	2		
5	C	1	Total	C	N	O	0	0
			8	5	1	2		
5	D	1	Total	C	N	O	0	0
			8	5	1	2		

- Molecule 6 is ALANINE (CCD ID: ALA) (formula: $C_3H_7NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			6	3	1	2		
6	B	1	Total	C	N	O	0	0
			5	3	1	1		

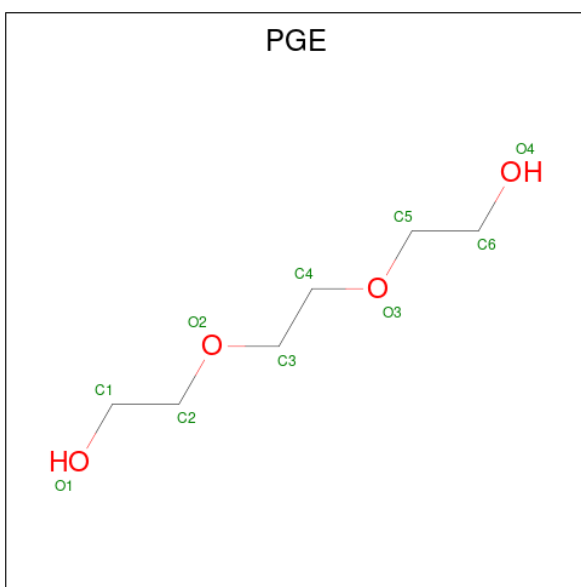
- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Ca	0	0
			1	1		
7	B	2	Total	Ca	0	0
			2	2		
7	C	2	Total	Ca	0	0
			2	2		

- Molecule 8 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

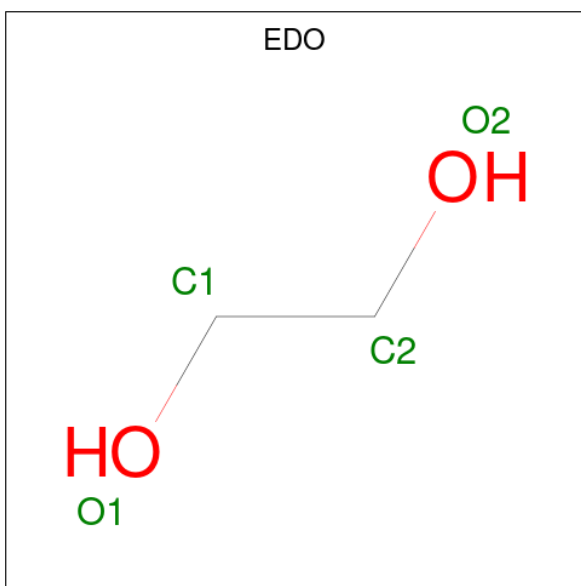
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	3	Total	Mn	0	0
			3	3		
8	B	3	Total	Mn	0	0
			3	3		
8	C	3	Total	Mn	0	0
			3	3		
8	D	3	Total	Mn	0	0
			3	3		

- Molecule 9 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 10 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	C	1	Total	C	O	0	0
			4	2	2		
10	D	1	Total	C	O	0	0
			4	2	2		

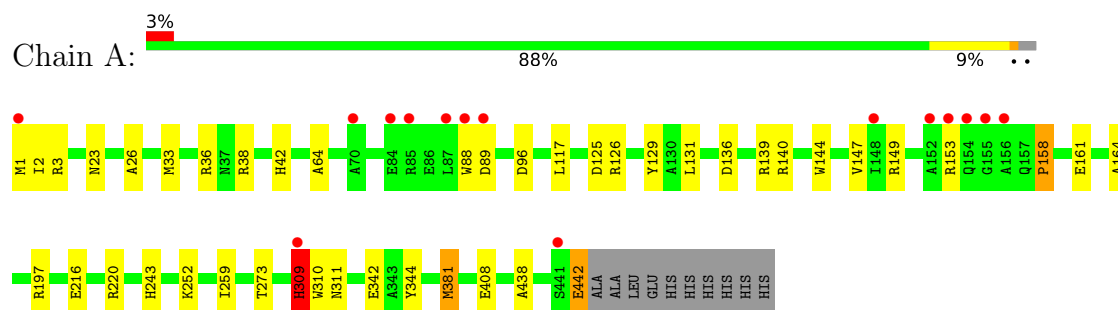
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	303	Total	O	0	0
			303	303		
11	B	275	Total	O	0	0
			275	275		
11	C	268	Total	O	0	0
			268	268		
11	D	306	Total	O	0	0
			306	306		

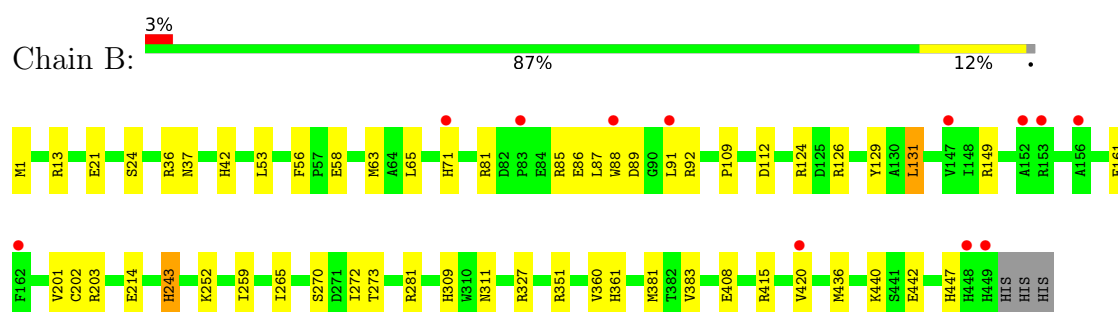
3 Residue-property plots [i](#)

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.

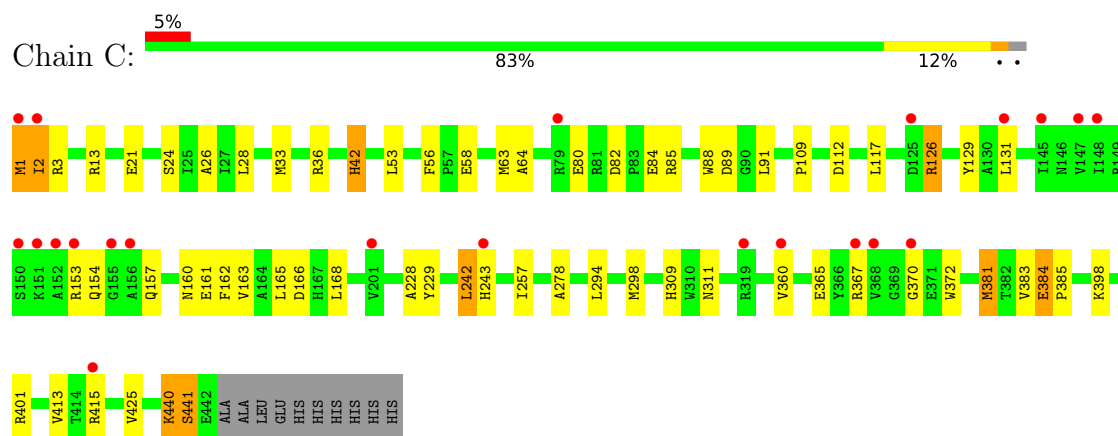
• Molecule 1: Aminopeptidase P



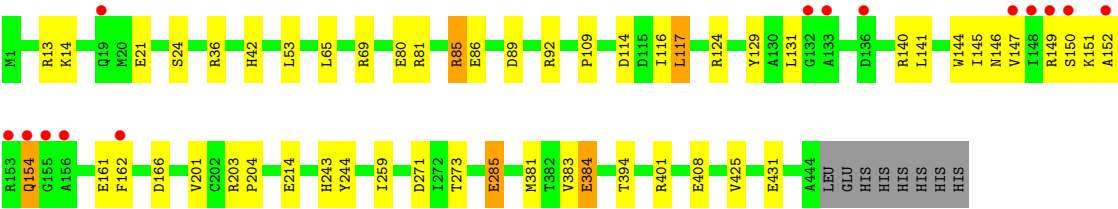
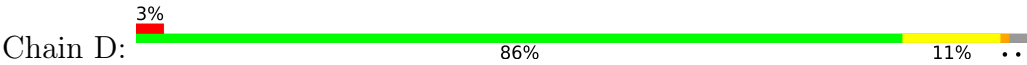
• Molecule 1: Aminopeptidase P



• Molecule 1: Aminopeptidase P



• Molecule 1: Aminopeptidase P



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	111.20Å 123.43Å 149.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.01 – 1.78 48.01 – 1.78	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.01-1.78) 99.4 (48.01-1.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 1.78Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.210 , 0.221 (Not available) , 0.211	Depositor DCC
R_{free} test set	2000 reflections (1.02%)	wwPDB-VP
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.229	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15378	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.85% 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: GOL, EDO, NA, PGE, CA, SO4, MN

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	0.62	5/3603 (0.1%)	0.66	3/4880 (0.1%)
1	B	0.56	2/3642 (0.1%)	0.57	4/4937 (0.1%)
1	C	0.56	3/3624 (0.1%)	0.60	4/4911 (0.1%)
1	D	0.63	7/3597 (0.2%)	0.64	4/4875 (0.1%)
All	All	0.59	17/14466 (0.1%)	0.62	15/19603 (0.1%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	309[A]	HIS	CA-C	7.69	1.64	1.53
1	A	309[B]	HIS	CA-C	7.69	1.64	1.53
1	A	117	LEU	C-O	-7.54	1.17	1.24
1	D	285	GLU	CA-C	6.50	1.61	1.52
1	D	285	GLU	N-CA	6.47	1.54	1.46
1	C	117	LEU	C-O	-6.11	1.18	1.24
1	B	281	ARG	C-O	-5.69	1.17	1.23
1	A	158	PRO	C-O	-5.61	1.18	1.24
1	D	384	GLU	C-O	-5.47	1.17	1.24
1	D	117	LEU	C-O	-5.39	1.19	1.24
1	C	385	PRO	C-O	-5.32	1.17	1.23
1	C	383	VAL	C-O	-5.27	1.18	1.24
1	D	244	TYR	C-O	-5.21	1.17	1.23
1	D	271	ASP	C-O	-5.21	1.17	1.23
1	D	383	VAL	C-O	-5.19	1.18	1.24
1	A	96	ASP	C-O	-5.15	1.18	1.24
1	B	243	HIS	C-O	-5.08	1.17	1.24

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	89	ASP	N-CA-C	12.86	125.38	111.36
1	D	89	ASP	N-CA-C	8.54	123.88	113.38
1	D	116	ILE	N-CA-C	8.10	118.92	112.12
1	A	309[A]	HIS	CA-C-O	6.63	129.54	121.72
1	A	309[B]	HIS	CA-C-O	6.63	129.54	121.72
1	C	440	LYS	N-CA-C	6.42	118.81	111.11
1	C	89	ASP	N-CA-C	6.06	120.67	113.28
1	C	384	GLU	N-CA-C	5.88	118.06	109.30
1	B	87	LEU	N-CA-C	-5.80	104.96	111.28
1	C	242	LEU	N-CA-C	5.58	117.36	111.28
1	B	89	ASP	N-CA-C	5.26	119.85	113.38
1	D	285	GLU	N-CA-C	5.23	117.06	111.36
1	D	285	GLU	CA-C-O	5.12	125.85	120.42
1	B	270	SER	N-CA-C	-5.11	101.43	109.50
1	B	272	ILE	N-CA-C	5.06	115.77	108.48

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	3511	0	3442	37	1
1	B	3550	0	3478	41	0
1	C	3523	0	3471	53	1
1	D	3509	0	3441	31	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
3	A	6	0	8	2	0
3	D	18	0	24	4	0
4	A	5	0	0	1	0
4	D	10	0	0	0	0
5	A	8	0	7	3	0
5	B	8	0	7	1	0
5	C	8	0	7	1	0
5	D	8	0	7	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	6	0	4	0	0
6	B	5	0	4	1	0
7	A	1	0	0	0	0
7	B	2	0	0	0	0
7	C	2	0	0	0	0
8	A	3	0	0	0	0
8	B	3	0	0	0	0
8	C	3	0	0	0	0
8	D	3	0	0	0	0
9	B	10	0	14	5	0
10	B	8	0	12	7	0
10	C	8	0	12	2	0
10	D	4	0	6	0	0
11	A	303	0	0	7	0
11	B	275	0	0	4	0
11	C	268	0	0	17	0
11	D	306	0	0	4	0
All	All	15378	0	13944	159	1

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 (159) 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:C:33:MET:HE3	11:C:826:HOH:O	1.51	1.10
1:B:243:HIS:NE2	5:B:506:PRO:HD2	1.82	0.95
1:A:38:ARG:HH12	3:A:502:GOL:H31	1.30	0.93
1:C:160:ASN:HB3	11:C:818:HOH:O	1.71	0.90
1:C:440:LYS:HE2	11:C:798:HOH:O	1.76	0.85
1:A:309[A]:HIS:HD2	1:A:311:ASN:H	1.25	0.84
1:D:201[B]:VAL:HG11	1:D:214:GLU:HG3	1.60	0.83
5:A:504:PRO:OXT	1:C:88:TRP:HZ3	1.61	0.81
1:A:33[B]:MET:HE3	11:A:781:HOH:O	1.80	0.80
1:C:367:ARG:NH2	1:C:370:GLY:O	2.15	0.80
1:B:327:ARG:NH1	11:B:601:HOH:O	2.14	0.80
1:B:203:ARG:HB3	9:B:503:PGE:H3	1.62	0.79
1:C:28[B]:LEU:HD22	11:C:856:HOH:O	1.81	0.79
1:A:438:ALA:O	1:A:442:GLU:HG2	1.82	0.79
1:C:309[A]:HIS:HE1	1:C:311:ASN:HB3	1.49	0.78
1:A:33[B]:MET:CE	11:A:781:HOH:O	2.37	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:HIS:CE1	10:B:505:EDO:H21	2.25	0.71
1:C:154:GLN:OE1	1:C:154:GLN:N	2.24	0.71
1:C:33:MET:CE	11:C:826:HOH:O	2.18	0.70
1:A:140[B]:ARG:NH2	11:A:602:HOH:O	2.24	0.70
1:B:252:LYS:HE2	11:B:621:HOH:O	1.92	0.69
1:A:38:ARG:NH1	3:A:502:GOL:H31	2.07	0.68
1:C:82:ASP:OD2	1:C:85:ARG:HD3	1.95	0.67
1:A:136:ASP:OD1	1:A:139:ARG:NH1	2.26	0.67
1:B:201[A]:VAL:HG11	1:B:214:GLU:HG3	1.76	0.67
1:B:1:MET:HE2	11:B:604:HOH:O	1.94	0.67
1:C:2:ILE:HD12	1:C:3:ARG:H	1.63	0.64
1:B:129:TYR:CZ	1:B:131:LEU:HD13	2.32	0.64
1:D:203:ARG:NH2	11:D:605:HOH:O	2.30	0.63
1:C:33:MET:HB3	11:C:826:HOH:O	1.98	0.63
5:A:504:PRO:OXT	1:C:88:TRP:CZ3	2.50	0.62
1:C:28[A]:LEU:HD12	11:C:856:HOH:O	1.99	0.62
1:C:372:TRP:CD1	10:C:503:EDO:H21	2.34	0.62
1:A:309[A]:HIS:CD2	1:A:311:ASN:H	2.13	0.61
1:A:88:TRP:O	1:C:243:HIS:HB2	2.01	0.60
1:B:126:ARG:HG2	1:B:161:GLU:HB2	1.84	0.60
1:B:1:MET:HE3	1:B:265:ILE:HG23	1.84	0.60
1:D:203:ARG:NH1	1:D:204:PRO:O	2.35	0.60
1:D:36:ARG:HG3	1:D:42:HIS:CE1	2.37	0.59
1:C:229:TYR:HD1	1:C:360[A]:VAL:HG11	1.69	0.58
1:D:69:ARG:HA	3:D:501:GOL:H11	1.85	0.58
1:C:33:MET:CG	11:C:826:HOH:O	2.52	0.58
1:A:126:ARG:HG2	1:A:161:GLU:HB2	1.86	0.57
1:C:28[B]:LEU:HD21	1:C:165:LEU:HD23	1.86	0.57
1:B:85:ARG:HG2	1:B:92:ARG:NH2	2.20	0.57
1:C:381:MET:HE1	1:C:413:VAL:HG23	1.88	0.56
1:D:151:LYS:O	1:D:152:ALA:C	2.48	0.56
1:C:360[B]:VAL:HG23	11:C:777:HOH:O	2.05	0.56
1:C:441:SER:OG	11:C:601:HOH:O	2.00	0.56
1:B:21:GLU:O	1:B:24:SER:OG	2.21	0.56
1:B:21:GLU:HG3	1:B:126:ARG:NH1	2.21	0.56
1:B:436:MET:HB3	9:B:503:PGE:H62	1.87	0.56
1:A:252:LYS:HE2	11:A:611:HOH:O	2.06	0.56
1:B:361:HIS:HE1	10:B:505:EDO:H22	1.71	0.56
1:A:220:ARG:HD3	11:C:729:HOH:O	2.06	0.56
1:C:21:GLU:O	1:C:24:SER:OG	2.21	0.56
1:C:80:GLU:HG3	1:C:109:PRO:HB3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:285:GLU:H	1:D:285:GLU:CD	2.14	0.55
1:D:401:ARG:HH22	3:D:503:GOL:H2	1.72	0.54
1:B:243:HIS:HE1	10:B:505:EDO:H21	1.70	0.54
1:C:126:ARG:HD3	1:C:163:VAL:HG21	1.88	0.54
1:C:242:LEU:O	11:C:602:HOH:O	2.18	0.54
1:A:309[A]:HIS:HD2	1:A:311:ASN:N	2.01	0.54
1:B:81:ARG:NH1	1:B:86:GLU:OE1	2.42	0.53
1:D:85:ARG:HD3	1:D:92:ARG:NH2	2.23	0.53
1:B:415:ARG:HG2	1:B:415:ARG:HH21	1.72	0.53
1:C:33:MET:HG2	11:C:826:HOH:O	2.08	0.53
1:D:114:ASP:CG	1:D:140:ARG:HH22	2.17	0.52
1:C:309[A]:HIS:CE1	1:C:311:ASN:HB3	2.37	0.52
1:A:36:ARG:HG3	1:A:42:HIS:CE1	2.44	0.52
1:B:351:ARG:HG3	6:B:507:ALA:HB3	1.92	0.52
1:A:216:GLU:O	1:A:220:ARG:HG3	2.10	0.52
1:A:129:TYR:CZ	1:A:131:LEU:HD13	2.45	0.51
1:B:36:ARG:HD2	1:B:37:ASN:OD1	2.11	0.51
1:C:28[A]:LEU:HB2	1:C:63:MET:HG2	1.92	0.51
1:B:13:ARG:HG2	1:B:65:LEU:HD21	1.92	0.51
1:D:166:ASP:HB2	11:D:692:HOH:O	2.10	0.50
1:A:243:HIS:HE2	5:A:504:PRO:N	2.07	0.50
1:D:151:LYS:O	1:D:154:GLN:N	2.44	0.50
1:A:131:LEU:HD23	1:A:164:ALA:HA	1.93	0.50
1:B:383:VAL:O	1:B:408:GLU:HA	2.11	0.50
1:C:28[A]:LEU:HD11	1:C:165:LEU:HD23	1.93	0.49
1:C:398:LYS:HG2	1:C:401:ARG:CZ	2.43	0.49
1:B:243:HIS:HE1	10:B:505:EDO:C2	2.25	0.48
1:C:84:GLU:HB2	11:C:704:HOH:O	2.12	0.48
1:C:372:TRP:HD1	10:C:503:EDO:H21	1.78	0.48
1:D:147:VAL:O	1:D:150:SER:OG	2.30	0.48
1:C:56:PHE:CE2	1:C:58:GLU:HB2	2.48	0.48
1:B:124:ARG:HH22	10:B:504:EDO:H22	1.77	0.48
1:D:431:GLU:OE1	11:D:601:HOH:O	2.19	0.48
1:A:3:ARG:NH2	4:A:503:SO4:O1	2.47	0.47
1:B:124:ARG:NH1	11:B:610:HOH:O	2.47	0.47
1:B:71:HIS:O	1:B:71:HIS:ND1	2.43	0.47
1:C:80:GLU:CG	1:C:109:PRO:HB3	2.44	0.47
1:A:88:TRP:O	1:C:243:HIS:N	2.45	0.47
1:D:124:ARG:HH22	3:D:501:GOL:H31	1.80	0.47
1:B:309:HIS:HE1	1:B:311:ASN:HB3	1.80	0.47
1:C:36:ARG:HG3	1:C:42:HIS:NE2	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:13:ARG:HD3	1:D:53:LEU:O	2.14	0.47
1:D:21:GLU:O	1:D:24:SER:OG	2.28	0.46
1:B:88:TRP:O	1:D:243:HIS:HB2	2.14	0.46
1:C:160:ASN:OD1	1:C:160:ASN:N	2.41	0.46
1:D:141:LEU:HA	1:D:144:TRP:CE3	2.51	0.46
1:A:144:TRP:O	1:A:147:VAL:HG22	2.16	0.46
1:A:273:THR:H	1:A:408:GLU:HB3	1.81	0.46
1:C:21:GLU:HG3	1:C:126:ARG:CZ	2.46	0.45
1:C:228:ALA:HB1	1:C:360[A]:VAL:CG1	2.47	0.45
1:D:161:GLU:C	1:D:162:PHE:HD1	2.25	0.45
1:B:243:HIS:CE1	10:B:505:EDO:C2	2.98	0.44
1:D:384:GLU:HB3	1:D:408:GLU:HB2	1.98	0.44
1:C:129:TYR:CZ	1:C:131:LEU:HG	2.52	0.44
1:B:124:ARG:HH22	10:B:504:EDO:C2	2.30	0.44
1:D:401:ARG:HH12	3:D:503:GOL:H2	1.82	0.44
1:B:63:MET:HE2	1:B:65:LEU:HG	2.00	0.44
1:B:13:ARG:HD3	1:B:53:LEU:O	2.18	0.44
1:B:36:ARG:HG3	1:B:42:HIS:CE1	2.53	0.44
1:A:252:LYS:CE	11:A:611:HOH:O	2.65	0.43
1:D:146:ASN:HA	1:D:149:ARG:HG3	2.00	0.43
1:C:28[B]:LEU:CD2	11:C:856:HOH:O	2.53	0.43
1:B:109:PRO:HG2	1:B:112:ASP:OD2	2.18	0.43
1:D:114:ASP:OD1	1:D:140:ARG:NH2	2.47	0.43
1:B:440:LYS:HG3	9:B:503:PGE:H22	2.00	0.43
1:A:149:ARG:HG2	1:A:158:PRO:HG3	2.00	0.43
1:B:440:LYS:HB2	9:B:503:PGE:H52	2.01	0.43
1:C:384:GLU:OE1	5:C:504:PRO:HA	2.19	0.42
1:A:252:LYS:NZ	11:A:611:HOH:O	2.47	0.42
1:C:13:ARG:HD3	1:C:53:LEU:O	2.19	0.42
1:D:13:ARG:HG2	1:D:65:LEU:HD21	2.00	0.42
1:C:1:MET:HB3	1:C:1:MET:HE3	1.60	0.42
1:A:342:GLU:HA	1:A:344:TYR:CE1	2.54	0.42
1:D:14:LYS:NZ	11:D:606:HOH:O	2.32	0.42
1:C:257:ILE:CD1	1:C:278:ALA:HB2	2.51	0.41
1:D:117:LEU:HD23	1:D:144:TRP:CZ3	2.55	0.41
1:A:1:MET:HG2	1:A:2:ILE:H	1.84	0.41
1:C:398:LYS:HG3	11:C:771:HOH:O	2.19	0.41
1:D:129:TYR:CZ	1:D:131:LEU:HG	2.56	0.41
1:C:228:ALA:HB1	1:C:360[B]:VAL:HG22	2.02	0.41
1:D:81:ARG:HD2	1:D:86:GLU:OE2	2.20	0.41
1:B:202:CYS:O	9:B:503:PGE:O4	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140[B]:ARG:HD2	1:A:144:TRP:CZ2	2.56	0.41
1:A:381:MET:HE2	11:A:822:HOH:O	2.20	0.41
1:A:259:ILE:O	1:A:273:THR:HA	2.21	0.41
1:B:1:MET:CE	1:B:265:ILE:HG23	2.49	0.41
1:B:56:PHE:CE2	1:B:58:GLU:HB2	2.55	0.41
1:C:26:ALA:O	1:C:64:ALA:HA	2.21	0.41
1:C:109:PRO:HG2	1:C:112:ASP:OD2	2.21	0.41
1:C:161:GLU:O	1:C:162:PHE:HD1	2.04	0.41
1:A:23:ASN:ND2	1:A:125:ASP:OD1	2.50	0.40
1:A:197:ARG:HD2	1:A:197:ARG:HA	1.94	0.40
1:C:298:MET:CE	11:C:865:HOH:O	2.69	0.40
1:D:80:GLU:HG3	1:D:109:PRO:HB3	2.03	0.40
1:A:26:ALA:O	1:A:64:ALA:HA	2.22	0.40
1:A:140[A]:ARG:HG2	1:A:144:TRP:CZ2	2.56	0.40
1:B:259:ILE:O	1:B:273:THR:HA	2.21	0.40
1:C:294:LEU:HD11	1:C:298:MET:HE1	2.04	0.40
1:D:259:ILE:O	1:D:273:THR:HA	2.21	0.40
1:B:442:GLU:OE1	1:B:447:HIS:ND1	2.49	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:ASP:OD2	1:C:415:ARG:NH2[2_354]	2.03	0.17

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	444/452 (98%)	436 (98%)	8 (2%)	0	100	100
1	B	451/452 (100%)	445 (99%)	6 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	447/452 (99%)	441 (99%)	6 (1%)	0	100	100
1	D	445/452 (98%)	438 (98%)	7 (2%)	0	100	100
All	All	1787/1808 (99%)	1760 (98%)	27 (2%)	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	355/359 (99%)	350 (99%)	5 (1%)	59	44
1	B	358/359 (100%)	350 (98%)	8 (2%)	45	26
1	C	357/359 (99%)	344 (96%)	13 (4%)	31	10
1	D	354/359 (99%)	348 (98%)	6 (2%)	53	36
All	All	1424/1436 (99%)	1392 (98%)	32 (2%)	47	26

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	ARG
1	A	309[A]	HIS
1	A	309[B]	HIS
1	A	381	MET
1	A	442	GLU
1	B	91	LEU
1	B	131	LEU
1	B	149	ARG
1	B	360[A]	VAL
1	B	360[B]	VAL
1	B	381	MET
1	B	420[A]	VAL
1	B	420[B]	VAL
1	C	1	MET

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Mol	Chain	Res	Type
1	C	2	ILE
1	C	42	HIS
1	C	91	LEU
1	C	126	ARG
1	C	153	ARG
1	C	157	GLN
1	C	166	ASP
1	C	168	LEU
1	C	365	GLU
1	C	381	MET
1	C	425	VAL
1	C	441	SER
1	D	85	ARG
1	D	145	ILE
1	D	154	GLN
1	D	381	MET
1	D	394	THR
1	D	425	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	HIS
1	A	157	GLN
1	A	311	ASN
1	B	172	GLN
1	B	243	HIS
1	D	361	HIS
1	D	416	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 40 ligands modelled in this entry, 21 are monoatomic - leaving 19 for Mogul analysis.

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	SO4	D	505	-	4,4,4	0.33	0	6,6,6	0.27	0
9	PGE	B	503	2	9,9,9	0.28	0	8,8,8	0.93	0
3	GOL	D	502	-	5,5,5	0.49	0	5,5,5	0.77	0
10	EDO	D	506	-	3,3,3	0.27	0	2,2,2	0.39	0
10	EDO	B	504	-	3,3,3	0.50	0	2,2,2	0.39	0
4	SO4	D	504	-	4,4,4	0.30	0	6,6,6	0.25	0
6	ALA	B	507	-	3,4,5	0.70	0	2,4,6	0.77	0
3	GOL	A	502	-	5,5,5	0.67	0	5,5,5	0.51	0
3	GOL	D	503	-	5,5,5	0.33	0	5,5,5	0.51	0
10	EDO	C	503	-	3,3,3	0.43	0	2,2,2	0.32	0
5	PRO	D	510	-	8,8,8	1.88	1 (12%)	10,10,10	1.37	1 (10%)
6	ALA	A	505	-	5,5,5	1.39	1 (20%)	6,6,6	1.81	1 (16%)
5	PRO	C	504	-	8,8,8	1.50	1 (12%)	10,10,10	2.27	3 (30%)
5	PRO	A	504	-	8,8,8	1.37	1 (12%)	10,10,10	2.54	3 (30%)
3	GOL	D	501	-	5,5,5	0.34	0	5,5,5	0.69	0
10	EDO	C	502	-	3,3,3	0.35	0	2,2,2	0.81	0
10	EDO	B	505	-	3,3,3	0.43	0	2,2,2	0.31	0
5	PRO	B	506	-	8,8,8	1.10	1 (12%)	10,10,10	1.53	2 (20%)
4	SO4	A	503	-	4,4,4	0.25	0	6,6,6	0.19	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
9	PGE	B	503	2	-	4/7/7/7	-
10	EDO	C	503	-	-	0/1/1/1	-
3	GOL	D	502	-	-	2/4/4/4	-
5	PRO	D	510	-	-	0/4/11/11	0/1/1/1
6	ALA	B	507	-	-	1/1/2/4	-
10	EDO	B	505	-	-	1/1/1/1	-
3	GOL	A	502	-	-	4/4/4/4	-
6	ALA	A	505	-	-	3/4/4/4	-
3	GOL	D	503	-	-	0/4/4/4	-
10	EDO	D	506	-	-	0/1/1/1	-
5	PRO	B	506	-	-	3/4/11/11	0/1/1/1
5	PRO	C	504	-	-	0/4/11/11	0/1/1/1
5	PRO	A	504	-	-	2/4/11/11	0/1/1/1
10	EDO	B	504	-	-	1/1/1/1	-
3	GOL	D	501	-	-	0/4/4/4	-
10	EDO	C	502	-	-	1/1/1/1	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	510	PRO	OXT-C	-4.76	1.15	1.30
5	C	504	PRO	OXT-C	-3.59	1.19	1.30
5	A	504	PRO	OXT-C	-3.32	1.20	1.30
6	A	505	ALA	OXT-C	-2.88	1.21	1.30
5	B	506	PRO	OXT-C	-2.09	1.24	1.30

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	504	PRO	OXT-C-O	-6.47	109.41	124.08
5	C	504	PRO	OXT-C-O	-5.49	111.63	124.08
6	A	505	ALA	OXT-C-O	-3.82	115.41	124.08
5	A	504	PRO	OXT-C-CA	3.60	125.70	113.51
5	B	506	PRO	OXT-C-O	-3.31	116.56	124.08
5	D	510	PRO	OXT-C-O	-2.92	117.46	124.08
5	C	504	PRO	C-CA-N	2.62	117.08	106.84
5	C	504	PRO	OXT-C-CA	2.55	122.14	113.51
5	B	506	PRO	OXT-C-CA	2.33	121.41	113.51
5	A	504	PRO	CD-N-CA	2.09	112.61	107.22

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	GOL	O1-C1-C2-O2
3	A	502	GOL	O1-C1-C2-C3
3	D	502	GOL	O1-C1-C2-C3
6	B	507	ALA	O-C-CA-CB
3	A	502	GOL	O2-C2-C3-O3
3	D	502	GOL	O1-C1-C2-O2
9	B	503	PGE	O3-C5-C6-O4
3	A	502	GOL	C1-C2-C3-O3
5	A	504	PRO	OXT-C-CA-N
5	A	504	PRO	O-C-CA-N
9	B	503	PGE	C3-C4-O3-C5
9	B	503	PGE	O2-C3-C4-O3
6	A	505	ALA	O-C-CA-CB
6	A	505	ALA	OXT-C-CA-CB
10	B	505	EDO	O1-C1-C2-O2
6	A	505	ALA	OXT-C-CA-N
10	C	502	EDO	O1-C1-C2-O2
5	B	506	PRO	O-C-CA-CB
10	B	504	EDO	O1-C1-C2-O2
5	B	506	PRO	O-C-CA-N
5	B	506	PRO	OXT-C-CA-N
9	B	503	PGE	C6-C5-O3-C4

There are no ring outliers.

12 monomers are involved in 27 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	B	503	PGE	5	0
10	B	504	EDO	2	0
6	B	507	ALA	1	0
3	A	502	GOL	2	0
3	D	503	GOL	2	0
10	C	503	EDO	2	0
5	C	504	PRO	1	0
5	A	504	PRO	3	0
3	D	501	GOL	2	0
10	B	505	EDO	5	0
5	B	506	PRO	1	0
4	A	503	SO4	1	0

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 ⓘ

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	442/452 (97%)	-0.11	15 (3%)	48	55	12, 21, 49, 92	4 (0%)
1	B	449/452 (99%)	-0.01	12 (2%)	56	63	12, 24, 54, 86	4 (0%)
1	C	442/452 (97%)	0.04	22 (4%)	34	40	14, 23, 53, 82	7 (1%)
1	D	444/452 (98%)	-0.04	14 (3%)	50	57	12, 22, 54, 101	3 (0%)
All	All	1777/1808 (98%)	-0.03	63 (3%)	47	53	12, 22, 53, 101	18 (1%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	319[A]	ARG	8.1
1	B	420[A]	VAL	8.0
1	B	152	ALA	5.8
1	A	88	TRP	5.5
1	C	153	ARG	3.7
1	D	156	ALA	3.7
1	A	155	GLY	3.4
1	C	147	VAL	3.3
1	A	152	ALA	3.3
1	B	449	HIS	3.1
1	A	156	ALA	3.1
1	D	152	ALA	3.1
1	C	148	ILE	3.0
1	C	156	ALA	2.9
1	B	91	LEU	2.9
1	C	201[A]	VAL	2.8
1	A	87	LEU	2.8
1	B	448	HIS	2.8
1	D	155	GLY	2.8
1	D	153	ARG	2.7
1	D	147	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	1	MET	2.7
1	A	1	MET	2.6
1	B	88	TRP	2.6
1	A	153	ARG	2.6
1	B	153	ARG	2.6
1	C	79[A]	ARG	2.6
1	C	415	ARG	2.6
1	A	89	ASP	2.6
1	C	152	ALA	2.6
1	C	360[A]	VAL	2.6
1	C	368	VAL	2.6
1	D	150	SER	2.6
1	B	83	PRO	2.5
1	C	2	ILE	2.5
1	A	441	SER	2.5
1	C	370	GLY	2.5
1	C	155	GLY	2.4
1	C	367	ARG	2.4
1	D	132	GLY	2.3
1	A	85	ARG	2.3
1	D	162	PHE	2.3
1	D	148	ILE	2.3
1	D	133	ALA	2.3
1	C	131	LEU	2.3
1	A	148	ILE	2.3
1	B	156	ALA	2.2
1	B	162	PHE	2.2
1	A	70	ALA	2.2
1	D	154	GLN	2.2
1	A	154	GLN	2.1
1	D	19	GLN	2.1
1	C	151	LYS	2.1
1	B	147	VAL	2.1
1	C	243	HIS	2.1
1	A	84	GLU	2.1
1	A	309[A]	HIS	2.1
1	C	145	ILE	2.1
1	D	136	ASP	2.0
1	C	150	SER	2.0
1	B	71	HIS	2.0
1	D	149	ARG	2.0
1	C	125	ASP	2.0

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
5	PRO	A	504	8/8	0.55	0.18	27,30,43,43	0
9	PGE	B	503	10/10	0.61	0.27	17,27,34,36	10
5	PRO	B	506	8/8	0.63	0.17	30,37,48,50	0
5	PRO	D	510	8/8	0.67	0.17	28,38,50,53	0
5	PRO	C	504	8/8	0.70	0.15	28,38,50,53	0
3	GOL	D	503	6/6	0.71	0.14	42,44,46,47	0
10	EDO	B	505	4/4	0.74	0.19	40,44,46,48	0
10	EDO	C	503	4/4	0.74	0.15	44,47,47,51	0
10	EDO	B	504	4/4	0.75	0.12	45,45,46,47	0
3	GOL	D	501	6/6	0.76	0.12	40,42,45,47	0
10	EDO	C	502	4/4	0.78	0.11	33,42,45,49	0
6	ALA	A	505	6/6	0.78	0.16	40,46,47,48	0
3	GOL	A	502	6/6	0.79	0.12	31,36,39,50	0
7	CA	B	508	1/1	0.83	0.14	80,80,80,80	0
6	ALA	B	507	5/6	0.83	0.12	34,37,40,44	0
10	EDO	D	506	4/4	0.83	0.12	38,39,43,45	0
7	CA	A	506	1/1	0.87	0.15	64,64,64,64	0
7	CA	B	509	1/1	0.87	0.14	76,76,76,76	0
3	GOL	D	502	6/6	0.88	0.09	27,38,39,40	0
7	CA	C	505	1/1	0.89	0.11	56,56,56,56	0
2	NA	C	501	1/1	0.90	0.18	36,36,36,36	0
8	MN	A	507	1/1	0.90	0.08	28,28,28,28	1
7	CA	C	506	1/1	0.94	0.10	48,48,48,48	0
8	MN	B	510	1/1	0.94	0.06	28,28,28,28	1
4	SO4	A	503	5/5	0.95	0.09	24,26,31,33	5
2	NA	B	501	1/1	0.95	0.17	34,34,34,34	0
8	MN	C	508	1/1	0.95	0.05	28,28,28,28	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	MN	D	507	1/1	0.95	0.06	37,37,37,37	0
2	NA	A	501	1/1	0.95	0.06	32,32,32,32	0
8	MN	B	512	1/1	0.96	0.05	33,33,33,33	0
8	MN	A	509	1/1	0.96	0.06	31,31,31,31	0
8	MN	B	511	1/1	0.96	0.04	28,28,28,28	0
8	MN	A	508	1/1	0.97	0.04	28,28,28,28	0
8	MN	D	509	1/1	0.97	0.05	28,28,28,28	1
8	MN	C	507	1/1	0.97	0.06	37,37,37,37	0
4	SO4	D	505	5/5	0.97	0.06	23,26,30,32	5
2	NA	B	502	1/1	0.98	0.12	32,32,32,32	0
8	MN	D	508	1/1	0.98	0.03	28,28,28,28	0
4	SO4	D	504	5/5	0.98	0.04	19,21,23,27	0
8	MN	C	509	1/1	0.99	0.03	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.