



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 10, 2023 – 12:23 PM EDT

PDB ID : 4K1W  
Title : Crystal structure of the A314P mutant of mannonate dehydratase from novosphingobium aromaticivorans complexed with mg and d-mannonate  
Authors : Fedorov, A.A.; Fedorov, E.V.; Wichelecki, D.; Gerlt, J.A.; Almo, S.C.  
Deposited on : 2013-04-06  
Resolution : 1.65 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35.1

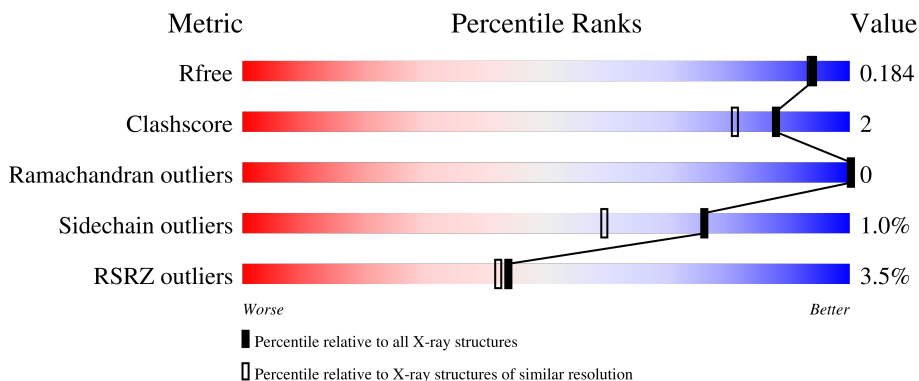
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.65 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))
$R_{free}$	130704	3122 (1.66-1.62)
Clashscore	141614	3268 (1.66-1.62)
Ramachandran outliers	138981	3215 (1.66-1.62)
Sidechain outliers	138945	3215 (1.66-1.62)
RSRZ outliers	127900	3079 (1.66-1.62)

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  $\geq 3$ , 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	418	 4% 89% 6% 6%
1	B	418	 2% 89% 8%
1	C	418	 3% 88% 8%
1	D	418	 4% 87% 8%

## 2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 13572 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 Mandelate racemase/muconate lactonizing enzyme, N-terminal domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	395	Total 3178	C 2032	N 552	O 577	S 17	0	7	0
1	B	386	Total 3092	C 1979	N 534	O 562	S 17	0	4	0
1	C	384	Total 3075	C 1968	N 535	O 555	S 17	0	2	0
1	D	384	Total 3088	C 1976	N 535	O 560	S 17	0	5	0

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	HIS	-	expression tag	UNP A4XF23
A	-14	HIS	-	expression tag	UNP A4XF23
A	-13	HIS	-	expression tag	UNP A4XF23
A	-12	HIS	-	expression tag	UNP A4XF23
A	-11	HIS	-	expression tag	UNP A4XF23
A	-10	HIS	-	expression tag	UNP A4XF23
A	-9	SER	-	expression tag	UNP A4XF23
A	-8	SER	-	expression tag	UNP A4XF23
A	-7	GLY	-	expression tag	UNP A4XF23
A	-6	LEU	-	expression tag	UNP A4XF23
A	-5	VAL	-	expression tag	UNP A4XF23
A	-4	PRO	-	expression tag	UNP A4XF23
A	-3	ARG	-	expression tag	UNP A4XF23
A	-2	GLY	-	expression tag	UNP A4XF23
A	-1	SER	-	expression tag	UNP A4XF23
A	0	HIS	-	expression tag	UNP A4XF23
A	314	PRO	ALA	engineered mutation	UNP A4XF23
B	-15	HIS	-	expression tag	UNP A4XF23
B	-14	HIS	-	expression tag	UNP A4XF23
B	-13	HIS	-	expression tag	UNP A4XF23

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	HIS	-	expression tag	UNP A4XF23
B	-11	HIS	-	expression tag	UNP A4XF23
B	-10	HIS	-	expression tag	UNP A4XF23
B	-9	SER	-	expression tag	UNP A4XF23
B	-8	SER	-	expression tag	UNP A4XF23
B	-7	GLY	-	expression tag	UNP A4XF23
B	-6	LEU	-	expression tag	UNP A4XF23
B	-5	VAL	-	expression tag	UNP A4XF23
B	-4	PRO	-	expression tag	UNP A4XF23
B	-3	ARG	-	expression tag	UNP A4XF23
B	-2	GLY	-	expression tag	UNP A4XF23
B	-1	SER	-	expression tag	UNP A4XF23
B	0	HIS	-	expression tag	UNP A4XF23
B	314	PRO	ALA	engineered mutation	UNP A4XF23
C	-15	HIS	-	expression tag	UNP A4XF23
C	-14	HIS	-	expression tag	UNP A4XF23
C	-13	HIS	-	expression tag	UNP A4XF23
C	-12	HIS	-	expression tag	UNP A4XF23
C	-11	HIS	-	expression tag	UNP A4XF23
C	-10	HIS	-	expression tag	UNP A4XF23
C	-9	SER	-	expression tag	UNP A4XF23
C	-8	SER	-	expression tag	UNP A4XF23
C	-7	GLY	-	expression tag	UNP A4XF23
C	-6	LEU	-	expression tag	UNP A4XF23
C	-5	VAL	-	expression tag	UNP A4XF23
C	-4	PRO	-	expression tag	UNP A4XF23
C	-3	ARG	-	expression tag	UNP A4XF23
C	-2	GLY	-	expression tag	UNP A4XF23
C	-1	SER	-	expression tag	UNP A4XF23
C	0	HIS	-	expression tag	UNP A4XF23
C	314	PRO	ALA	engineered mutation	UNP A4XF23
D	-15	HIS	-	expression tag	UNP A4XF23
D	-14	HIS	-	expression tag	UNP A4XF23
D	-13	HIS	-	expression tag	UNP A4XF23
D	-12	HIS	-	expression tag	UNP A4XF23
D	-11	HIS	-	expression tag	UNP A4XF23
D	-10	HIS	-	expression tag	UNP A4XF23
D	-9	SER	-	expression tag	UNP A4XF23
D	-8	SER	-	expression tag	UNP A4XF23
D	-7	GLY	-	expression tag	UNP A4XF23
D	-6	LEU	-	expression tag	UNP A4XF23
D	-5	VAL	-	expression tag	UNP A4XF23

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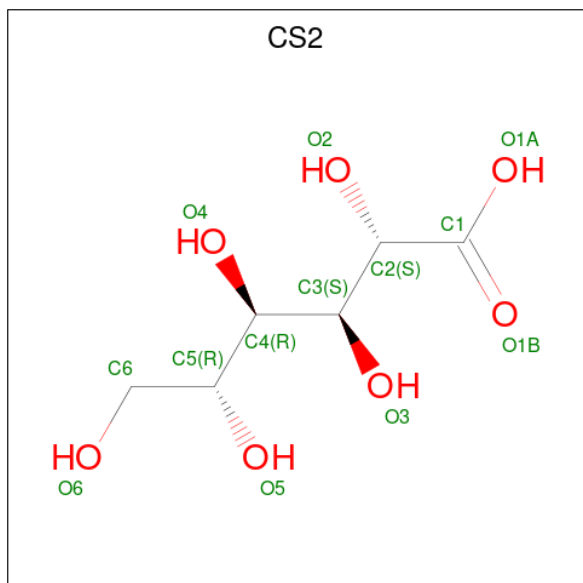
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Chain	Residue	Modelled	Actual	Comment	Reference
D	-4	PRO	-	expression tag	UNP A4XF23
D	-3	ARG	-	expression tag	UNP A4XF23
D	-2	GLY	-	expression tag	UNP A4XF23
D	-1	SER	-	expression tag	UNP A4XF23
D	0	HIS	-	expression tag	UNP A4XF23
D	314	PRO	ALA	engineered mutation	UNP A4XF23

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is D-MANNONIC ACID (three-letter code: CS2) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>7</sub>).



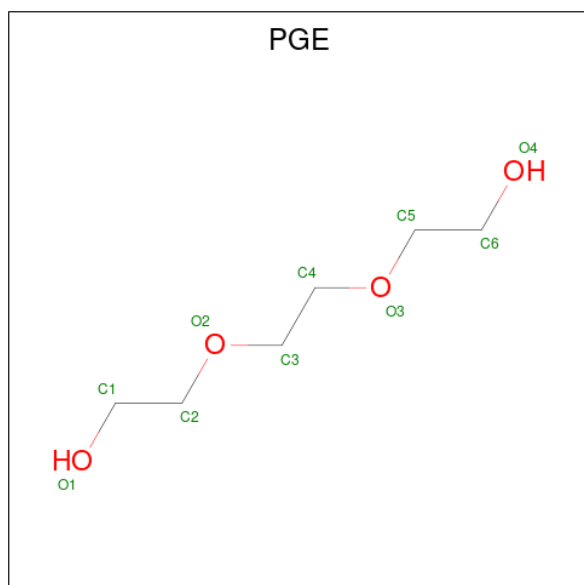
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 13 6 7	0	0
3	B	1	Total C O 13 6 7	0	0

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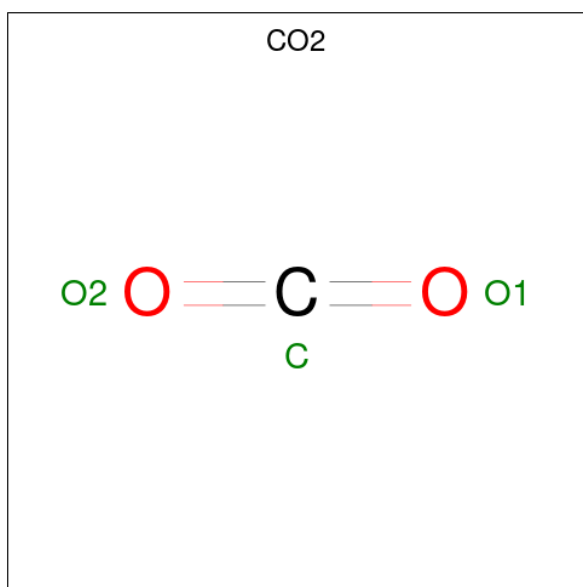
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			13	6	7		
3	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>4</sub>).



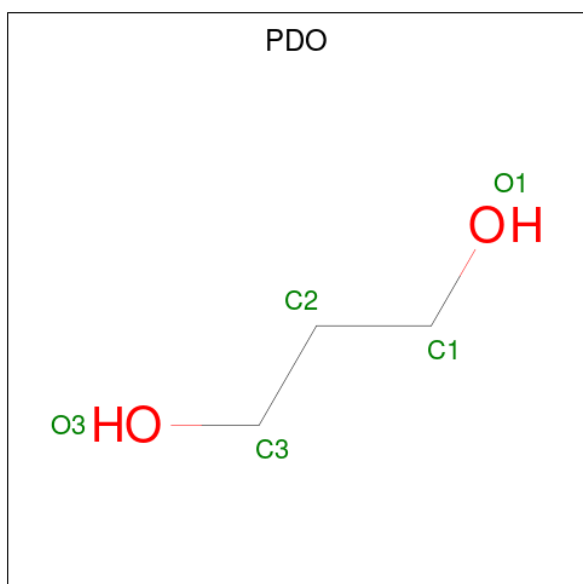
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		
4	C	1	Total	C	O	0	0
			10	6	4		
4	D	1	Total	C	O	0	0
			10	6	4		

- Molecule 5 is CARBON DIOXIDE (three-letter code: CO2) (formula: CO<sub>2</sub>).



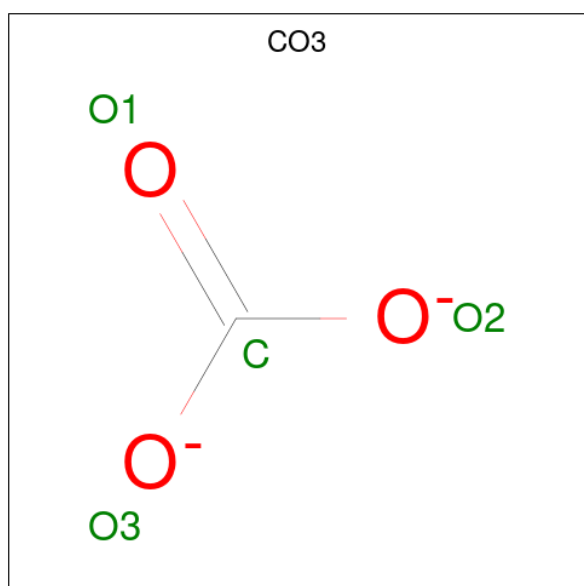
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 3 1 2	0	0
5	A	1	Total C O 3 1 2	0	0
5	C	1	Total C O 3 1 2	0	0
5	C	1	Total C O 3 1 2	0	0
5	D	1	Total C O 3 1 2	0	0

- Molecule 6 is 1,3-PROPANDIOL (three-letter code: PDO) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>2</sub>).



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

- Molecule 7 is CARBONATE ION (three-letter code: CO3) (formula: CO<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	C	1	Total	C	O	0	0
			4	1	3		

- Molecule 8 is water.

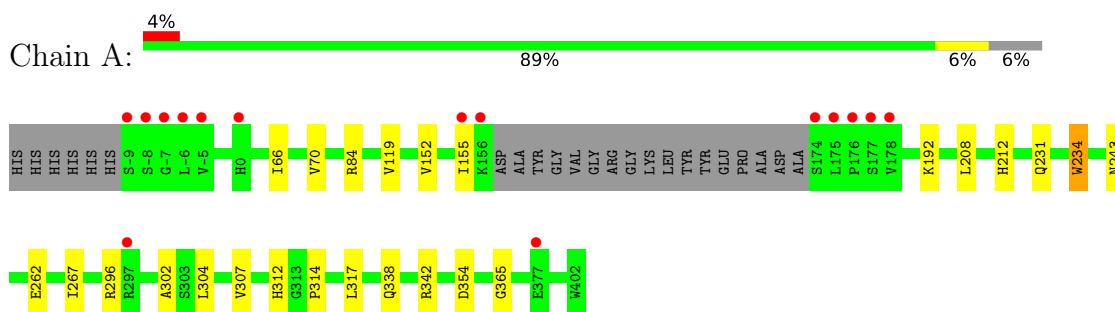
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	261	Total	O	0	0
			261	261		
8	B	288	Total	O	0	0
			288	288		
8	C	230	Total	O	0	0
			230	230		
8	D	224	Total	O	0	0
			224	224		



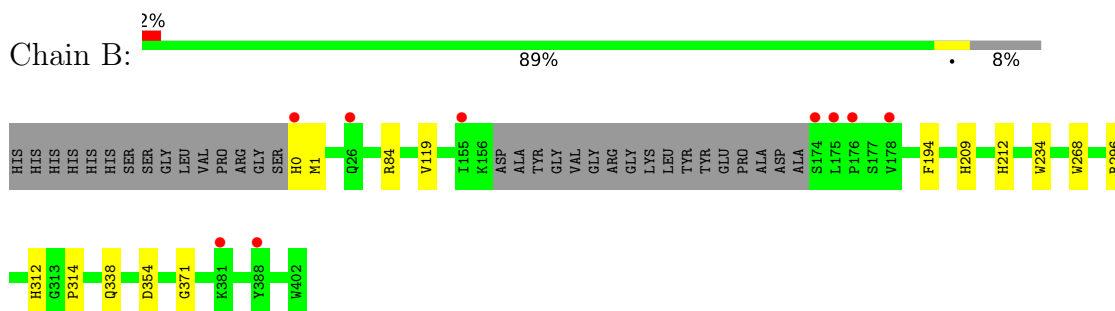
### 3 Residue-property plots [i](#)

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.

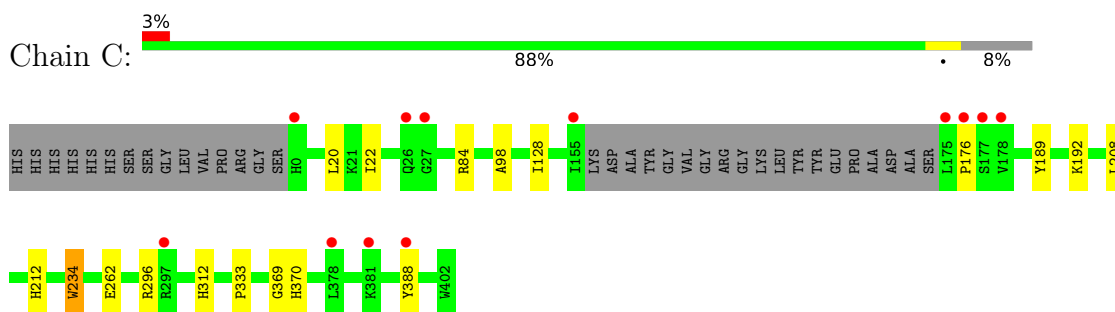
- Molecule 1: Mandelate racemase/muconate lactonizing enzyme, N-terminal domain protein



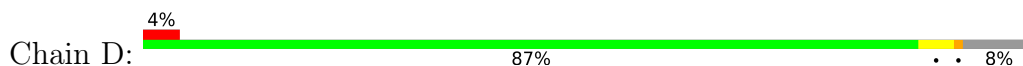
- Molecule 1: Mandelate racemase/muconate lactonizing enzyme, N-terminal domain protein

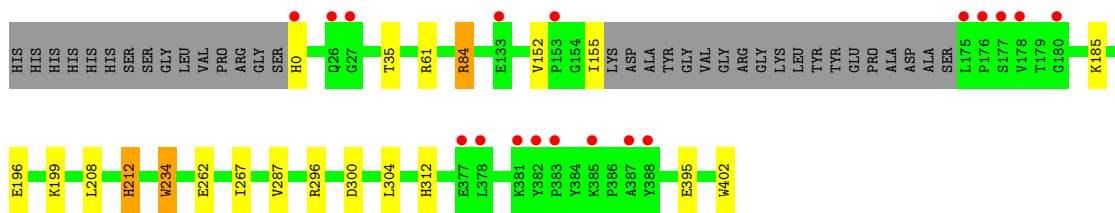


- Molecule 1: Mandelate racemase/muconate lactonizing enzyme, N-terminal domain protein



- Molecule 1: Mandelate racemase/muconate lactonizing enzyme, N-terminal domain protein





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	116.22Å 165.36Å 167.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.21 – 1.65 46.21 – 1.65	Depositor EDS
% Data completeness (in resolution range)	98.9 (46.21-1.65) 98.9 (46.21-1.65)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.44 (at 1.64Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, $R_{free}$	0.157 , 0.184 0.157 , 0.184	Depositor DCC
$R_{free}$ test set	9618 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.8	Xtrriage
Anisotropy	0.284	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 48.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	13572	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: PGE, PDO, CS2, CO2, CO3, MG

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	0.34	0/3276	0.54	0/4455
1	B	0.36	0/3189	0.55	0/4338
1	C	0.32	0/3163	0.51	0/4303
1	D	0.32	0/3185	0.53	0/4333
All	All	0.34	0/12813	0.53	0/17429

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	3178	0	3112	18	0
1	B	3092	0	3019	8	0
1	C	3075	0	3000	12	0
1	D	3088	0	3013	14	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	13	0	10	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	13	0	10	0	0
3	C	13	0	10	0	0
3	D	13	0	10	0	0
4	A	10	0	14	2	0
4	B	10	0	14	0	0
4	C	10	0	14	3	0
4	D	10	0	14	1	0
5	A	6	0	0	1	0
5	C	6	0	0	1	0
5	D	3	0	0	0	0
6	A	5	0	8	0	0
6	B	5	0	8	0	0
6	C	5	0	8	0	0
6	D	5	0	8	0	0
7	C	4	0	0	0	0
8	A	261	0	0	2	0
8	B	288	0	0	2	0
8	C	230	0	0	2	0
8	D	224	0	0	3	0
All	All	13572	0	12272	50	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 (50) 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:296:ARG:NH2	8:B:752:HOH:O	2.23	0.71
1:C:296:ARG:NH1	8:C:816:HOH:O	2.25	0.70
1:A:296:ARG:NH2	8:A:782:HOH:O	2.25	0.69
1:D:296:ARG:NH1	8:D:760:HOH:O	2.30	0.65
1:B:296:ARG:NH1	8:B:712:HOH:O	2.32	0.62
1:C:192:LYS:HZ3	4:C:503:PGE:H2	1.64	0.61
1:C:296:ARG:NH2	8:C:778:HOH:O	2.33	0.61
1:D:84:ARG:HD3	1:D:287:VAL:HG22	1.83	0.59
1:B:354[B]:ASP:OD1	1:B:371:GLY:HA3	2.08	0.53
1:D:296:ARG:NH2	8:D:772:HOH:O	2.43	0.52
1:A:192:LYS:HZ3	4:A:503:PGE:H2	1.76	0.51
1:D:212:HIS:ND1	1:D:402:TRP:OXT	2.36	0.51
1:C:192:LYS:HZ3	4:C:503:PGE:H42	1.77	0.49
1:A:192:LYS:HZ1	4:A:503:PGE:H6	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304[B]:LEU:HD21	1:D:267:ILE:HD13	1.95	0.48
1:A:342:ARG:NH2	8:A:807:HOH:O	2.48	0.47
1:D:196:GLU:HA	1:D:199:LYS:HG2	1.97	0.47
1:C:208:LEU:HD23	1:C:234:TRP:CE2	2.50	0.46
1:D:208:LEU:HD23	1:D:234:TRP:CE2	2.50	0.46
1:A:152:VAL:HB	1:A:155:ILE:HD12	1.99	0.45
1:D:300[B]:ASP:OD1	8:D:690:HOH:O	2.21	0.45
1:A:302:ALA:HB1	1:A:307:VAL:HB	1.99	0.44
1:A:314:PRO:HD2	1:A:317:LEU:HD22	1.99	0.44
1:C:128:ILE:HG13	1:C:189:TYR:CE2	2.53	0.44
1:D:185:LYS:NZ	1:D:395:GLU:OE1	2.39	0.44
1:D:35:THR:O	1:D:84:ARG:HD2	2.18	0.44
1:D:152:VAL:HB	1:D:155:ILE:HD12	2.00	0.44
1:A:208:LEU:HD23	1:A:234:TRP:CE2	2.52	0.43
1:C:192:LYS:NZ	4:C:503:PGE:H42	2.33	0.43
1:D:262:GLU:HB2	1:D:312:HIS:CG	2.53	0.43
1:C:20:LEU:HD21	1:C:22:ILE:HD11	2.00	0.43
1:C:333:PRO:HG2	5:C:505:CO2:C	2.49	0.42
1:B:0:HIS:ND1	1:B:1:MET:O	2.40	0.42
1:C:176:PRO:HD3	1:C:388:TYR:CG	2.54	0.42
1:D:0:HIS:HB2	1:D:61:ARG:NH2	2.34	0.42
1:A:312:HIS:CE1	1:A:314:PRO:HG3	2.54	0.42
1:C:98:ALA:HB3	1:C:369:GLY:HA2	2.02	0.42
4:D:503:PGE:H3	4:D:503:PGE:H1	1.83	0.42
1:A:119:VAL:HA	1:A:338:GLN:O	2.20	0.42
1:B:312:HIS:CE1	1:B:314:PRO:HG3	2.54	0.42
1:A:243:ASN:HB2	1:B:268:TRP:CZ3	2.55	0.41
1:A:262:GLU:HB2	1:A:312:HIS:CG	2.55	0.41
1:A:267:ILE:HD13	1:D:304:LEU:HD21	2.02	0.41
1:B:194:PHE:CZ	1:B:209:HIS:HB2	2.55	0.41
1:A:231:GLN:HE21	5:A:505:CO2:C	2.33	0.41
1:B:119:VAL:HA	1:B:338:GLN:O	2.21	0.41
1:A:66:ILE:O	1:A:70[A]:VAL:HG23	2.20	0.41
1:A:312:HIS:NE2	3:A:502:CS2:H3	2.35	0.41
1:A:354:ASP:CG	1:A:365:GLY:HA3	2.41	0.41
1:C:262:GLU:HB2	1:C:312:HIS:CG	2.56	0.41

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	398/418 (95%)	386 (97%)	12 (3%)	0	100	100
1	B	386/418 (92%)	373 (97%)	13 (3%)	0	100	100
1	C	382/418 (91%)	371 (97%)	11 (3%)	0	100	100
1	D	385/418 (92%)	373 (97%)	12 (3%)	0	100	100
All	All	1551/1672 (93%)	1503 (97%)	48 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	331/341 (97%)	328 (99%)	3 (1%)	78	63
1	B	321/341 (94%)	318 (99%)	3 (1%)	78	63
1	C	317/341 (93%)	313 (99%)	4 (1%)	69	47
1	D	320/341 (94%)	317 (99%)	3 (1%)	78	63
All	All	1289/1364 (94%)	1276 (99%)	13 (1%)	76	59

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	ARG
1	A	212	HIS

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Mol	Chain	Res	Type
1	A	234	TRP
1	B	84	ARG
1	B	212	HIS
1	B	234	TRP
1	C	84	ARG
1	C	212	HIS
1	C	234	TRP
1	C	370	HIS
1	D	84	ARG
1	D	212	HIS
1	D	234	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 23 ligands modelled in this entry, 5 are monoatomic - leaving 18 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
3	CS2	B	502	2	12,12,12	0.90	0	16,16,16	1.09	0
4	PGE	D	503	-	9,9,9	0.46	0	8,8,8	0.25	0
5	CO2	C	505	-	2,2,2	1.07	0	1,1,1	0.47	0
5	CO2	C	504	-	2,2,2	0.80	0	1,1,1	0.48	0
5	CO2	A	504	-	2,2,2	0.76	0	1,1,1	0.47	0
3	CS2	C	502	2	12,12,12	0.93	0	16,16,16	1.11	1 (6%)
3	CS2	A	502	2	12,12,12	1.08	1 (8%)	16,16,16	1.38	1 (6%)
5	CO2	D	504	-	2,2,2	1.02	0	1,1,1	0.38	0
6	PDO	A	506	-	4,4,4	0.56	0	3,3,3	0.72	0
4	PGE	B	503	-	9,9,9	0.47	0	8,8,8	0.29	0
5	CO2	A	505	-	2,2,2	1.03	0	1,1,1	0.51	0
4	PGE	C	503	-	9,9,9	0.48	0	8,8,8	0.24	0
6	PDO	D	505	-	4,4,4	0.46	0	3,3,3	0.99	0
3	CS2	D	502	2	12,12,12	0.81	0	16,16,16	1.41	2 (12%)
6	PDO	B	505	-	4,4,4	0.49	0	3,3,3	0.98	0
7	CO3	C	506	-	2,3,3	0.37	0	2,3,3	0.10	0
4	PGE	A	503	-	9,9,9	0.51	0	8,8,8	0.32	0
6	PDO	C	507	-	4,4,4	0.52	0	3,3,3	0.86	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
3	CS2	B	502	2	-	2/18/18/18	-
4	PGE	D	503	-	-	3/7/7/7	-
3	CS2	C	502	2	-	0/18/18/18	-
3	CS2	A	502	2	-	0/18/18/18	-
6	PDO	A	506	-	-	1/2/2/2	-
4	PGE	B	503	-	-	1/7/7/7	-
4	PGE	C	503	-	-	4/7/7/7	-
6	PDO	D	505	-	-	0/2/2/2	-
3	CS2	D	502	2	-	0/18/18/18	-
6	PDO	B	505	-	-	0/2/2/2	-
4	PGE	A	503	-	-	3/7/7/7	-
6	PDO	C	507	-	-	1/2/2/2	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	CS2	C5-C4	2.06	1.57	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	A	502	CS2	C4-C3-C2	4.14	120.94	113.60
3	D	502	CS2	C3-C2-C1	-3.32	105.15	109.32
3	C	502	CS2	C4-C3-C2	3.09	119.08	113.60
3	D	502	CS2	C4-C3-C2	2.30	117.68	113.60

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	503	PGE	C6-C5-O3-C4
4	D	503	PGE	O2-C3-C4-O3
4	A	503	PGE	O2-C3-C4-O3
4	A	503	PGE	C4-C3-O2-C2
4	A	503	PGE	C6-C5-O3-C4
4	B	503	PGE	O1-C1-C2-O2
4	D	503	PGE	C1-C2-O2-C3
4	C	503	PGE	O1-C1-C2-O2
4	D	503	PGE	C4-C3-O2-C2
3	B	502	CS2	O2-C2-C3-O3
3	B	502	CS2	O2-C2-C3-C4
4	C	503	PGE	C4-C3-O2-C2
4	C	503	PGE	O2-C3-C4-O3
6	A	506	PDO	C1-C2-C3-O3
6	C	507	PDO	O1-C1-C2-C3

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	503	PGE	1	0
5	C	505	CO2	1	0
3	A	502	CS2	1	0
5	A	505	CO2	1	0
4	C	503	PGE	3	0
4	A	503	PGE	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	395/418 (94%)	-0.22	15 (3%) 40 38	13, 24, 43, 66	0
1	B	386/418 (92%)	-0.32	9 (2%) 60 60	12, 21, 36, 63	0
1	C	384/418 (91%)	-0.02	12 (3%) 49 47	15, 30, 48, 69	0
1	D	384/418 (91%)	-0.01	18 (4%) 31 29	15, 28, 45, 68	0
All	All	1549/1672 (92%)	-0.15	54 (3%) 44 42	12, 25, 45, 69	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	175	LEU	6.9
1	C	175	LEU	5.9
1	A	-7	GLY	5.5
1	A	-8	SER	5.4
1	A	176	PRO	5.3
1	C	178	VAL	5.0
1	A	-6	LEU	5.0
1	D	0	HIS	5.0
1	D	177	SER	4.5
1	B	0	HIS	4.5
1	D	176	PRO	4.5
1	C	176	PRO	4.4
1	A	-5	VAL	4.4
1	D	385	LYS	4.3
1	A	175	LEU	4.2
1	A	178	VAL	4.2
1	A	-9	SER	4.2
1	C	0	HIS	4.1
1	B	175	LEU	3.9
1	D	382	TYR	3.6
1	B	174	SER	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	383	PRO	3.5
1	C	27	GLY	3.4
1	B	26	GLN	3.3
1	D	178	VAL	3.2
1	A	156	LYS	3.2
1	C	26	GLN	3.2
1	A	174	SER	3.1
1	A	177	SER	3.1
1	B	178	VAL	2.8
1	C	155	ILE	2.7
1	D	377	GLU	2.7
1	B	155	ILE	2.7
1	C	177	SER	2.6
1	D	153	PRO	2.6
1	C	297[A]	ARG	2.6
1	D	27	GLY	2.5
1	A	155	ILE	2.5
1	D	378	LEU	2.5
1	C	388	TYR	2.4
1	D	387	ALA	2.4
1	A	377	GLU	2.4
1	C	378	LEU	2.3
1	B	388	TYR	2.3
1	B	176	PRO	2.2
1	D	388	TYR	2.2
1	D	133	GLU	2.2
1	D	180	GLY	2.2
1	A	297[A]	ARG	2.1
1	D	381	LYS	2.1
1	A	0	HIS	2.1
1	D	26	GLN	2.1
1	C	381	LYS	2.0
1	B	381	LYS	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 monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	PGE	A	503	10/10	0.78	0.14	34,39,44,47	0
4	PGE	D	503	10/10	0.81	0.12	49,55,58,59	0
5	CO2	A	505	3/3	0.81	0.24	46,46,49,51	0
6	PDO	C	507	5/5	0.86	0.14	29,43,45,46	0
4	PGE	C	503	10/10	0.87	0.10	43,45,52,55	0
3	CS2	D	502	13/13	0.89	0.14	23,34,41,41	0
6	PDO	D	505	5/5	0.90	0.14	24,29,35,40	0
4	PGE	B	503	10/10	0.91	0.11	38,41,42,43	0
5	CO2	C	505	3/3	0.91	0.09	30,30,41,46	0
5	CO2	D	504	3/3	0.92	0.32	36,36,38,42	0
3	CS2	C	502	13/13	0.93	0.11	25,35,42,42	0
6	PDO	B	505	5/5	0.93	0.12	22,36,39,44	0
3	CS2	B	502	13/13	0.94	0.08	20,29,35,36	0
3	CS2	A	502	13/13	0.94	0.08	21,34,43,45	0
6	PDO	A	506	5/5	0.94	0.11	21,32,35,43	0
7	CO3	C	506	4/4	0.95	0.25	32,42,43,48	0
5	CO2	C	504	3/3	0.96	0.11	10,10,20,33	0
5	CO2	A	504	3/3	0.96	0.14	11,11,15,42	0
2	MG	A	501	1/1	0.98	0.06	19,19,19,19	0
2	MG	D	501	1/1	0.98	0.04	22,22,22,22	0
2	MG	B	504	1/1	0.99	0.03	16,16,16,16	1
2	MG	C	501	1/1	0.99	0.08	21,21,21,21	0
2	MG	B	501	1/1	0.99	0.04	18,18,18,18	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.