



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:18 AM GMT

PDB ID : 3OM6  
Title : Crystal structure of B. megaterium levansucrase mutant Y247A  
Authors : Strube, C.P.; Homann, A.; Gamer, M.; Jahn, D.; Seibel, J.; Heinz, D.W.  
Deposited on : 2010-08-26  
Resolution : 1.96 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

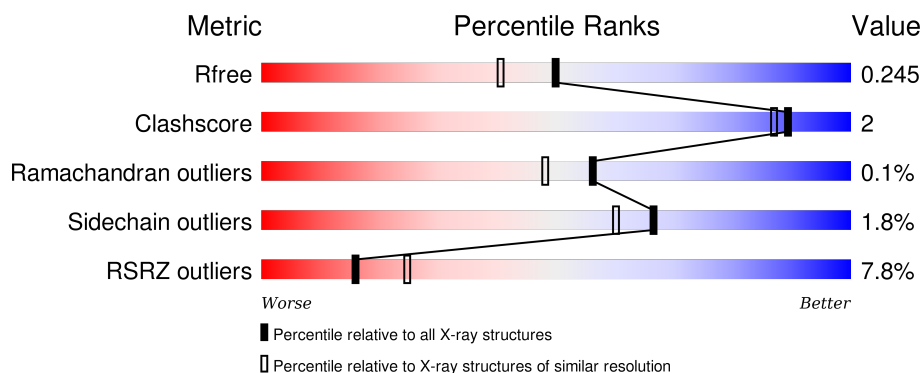
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.96 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))
$R_{free}$	91344	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	456	<div> <div>7%</div> <div>92%</div> <div>6%</div> </div>
1	B	456	<div> <div>10%</div> <div>89%</div> <div>9%</div> </div>
1	C	456	<div> <div>5%</div> <div>90%</div> <div>8%</div> </div>
1	D	456	<div> <div>9%</div> <div>91%</div> <div>7%</div> </div>

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
3	SO4	B	5	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 15500 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 Levansucrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	S	0	0	0
			3535	2218	592	718	7			
1	B	448	Total	C	N	O	S	0	2	0
			3552	2229	594	722	7			
1	C	448	Total	C	N	O	S	0	0	0
			3535	2218	592	718	7			
1	D	448	Total	C	N	O	S	0	4	0
			3562	2234	596	725	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	247	ALA	TYR	ENGINEERED MUTATION	UNP D5DC07
B	247	ALA	TYR	ENGINEERED MUTATION	UNP D5DC07
C	247	ALA	TYR	ENGINEERED MUTATION	UNP D5DC07
D	247	ALA	TYR	ENGINEERED MUTATION	UNP D5DC07

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

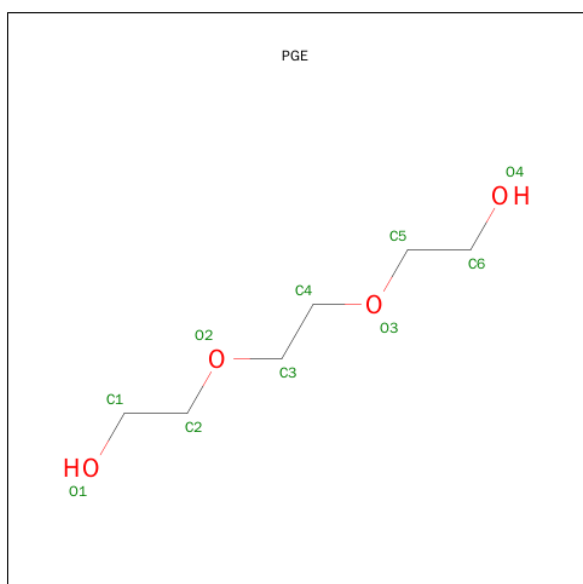
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		
2	D	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- 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	C	1	Total	C	O	0	0
			10	6	4		

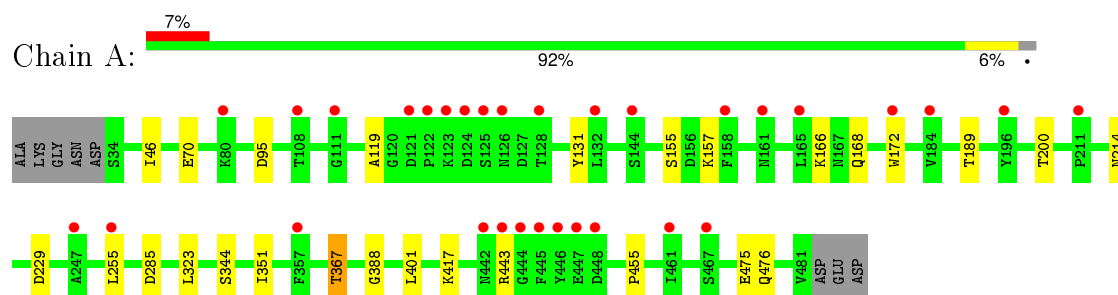
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	352	Total	O	0	0
			352	352		
5	B	347	Total	O	0	0
			347	347		
5	C	276	Total	O	0	0
			276	276		
5	D	302	Total	O	0	0
			302	302		

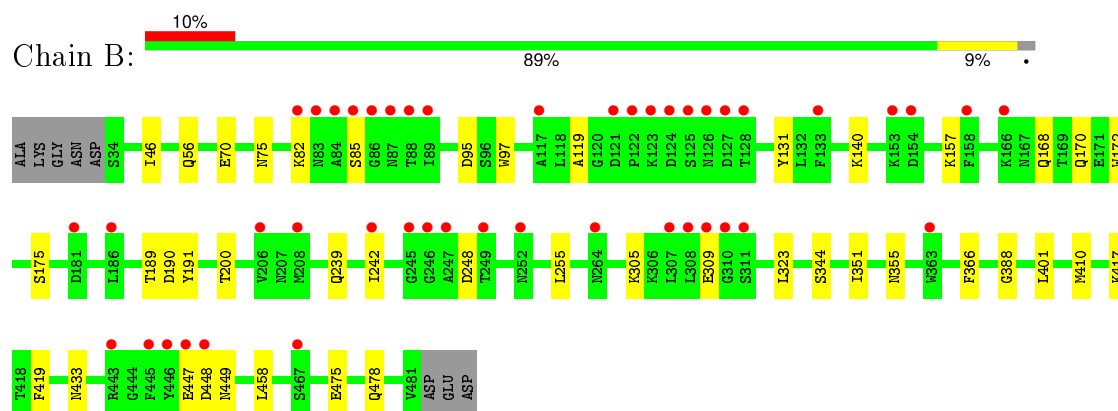
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

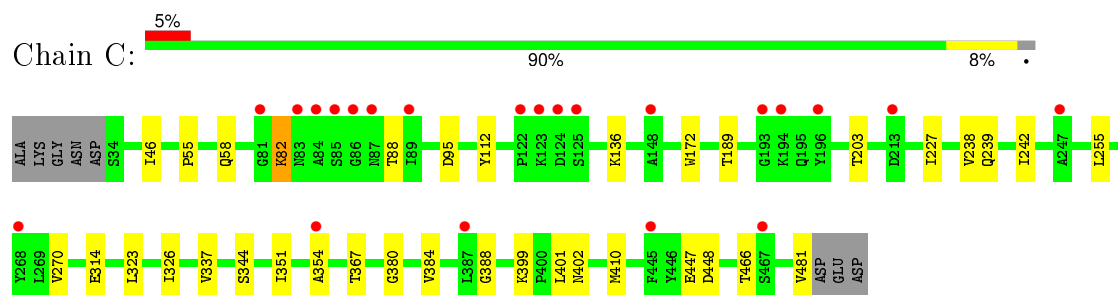
#### • Molecule 1: Levansucrase



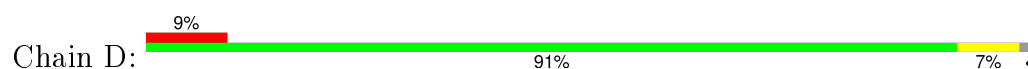
#### • Molecule 1: Levansucrase

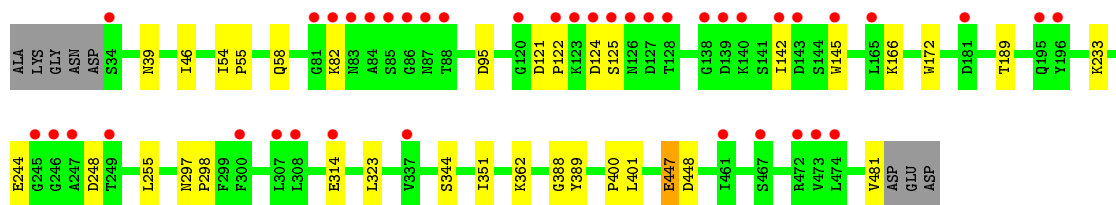


#### • Molecule 1: Levansucrase



#### • Molecule 1: Levansucrase







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.79Å 100.02Å 95.54Å 90.00° 90.59° 90.00°	Depositor
Resolution (Å)	50.01 – 1.96 47.77 – 1.96	Depositor EDS
% Data completeness (in resolution range)	93.5 (50.01-1.96) 95.2 (47.77-1.96)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.05 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.222 , 0.230 0.230 , 0.245	Depositor DCC
$R_{free}$ test set	6203 reflections (5.43%)	DCC
Wilson B-factor (Å <sup>2</sup> )	16.1	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 40.3	EDS
Estimated twinning fraction	0.664 for H, K, L 0.116 for -H, -K, L 0.110 for L, -K, H 0.111 for -L, K, H 0.012 for l,k,-h 0.028 for h,-k,-l 0.019 for l,-k,h	Xtriage
Reported twinning fraction	0.664 for H, K, L 0.116 for -H, -K, L 0.110 for L, -K, H 0.111 for -L, K, H	Depositor
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 120932 reflections	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	15500	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CA, PGE, SO4

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	0/3611	0.48	0/4887
1	B	0.34	0/3628	0.48	0/4910
1	C	0.34	0/3611	0.48	0/4887
1	D	0.35	0/3647	0.48	0/4935
All	All	0.34	0/14497	0.48	0/19619

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	3535	0	3374	13	0
1	B	3552	0	3389	21	0
1	C	3535	0	3374	20	0
1	D	3562	0	3400	15	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	15	0	0	0	0
3	D	5	0	0	0	0
4	C	10	0	14	0	0
5	A	352	0	0	1	0
5	B	347	0	0	0	0
5	C	276	0	0	1	0
5	D	302	0	0	1	0
All	All	15500	0	13551	67	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 (67) 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:305:LYS:O	1:B:309[A]:GLU:HG3	1.77	0.85
1:C:402:ASN:HD21	1:C:466:THR:H	1.26	0.82
1:A:155:SER:HB2	1:B:85:SER:HA	1.68	0.76
1:A:351:ILE:HB	1:A:367:THR:HG21	1.73	0.69
1:D:314[B]:GLU:OE1	1:D:314[B]:GLU:HA	1.93	0.68
1:C:82:LYS:HE3	1:C:82:LYS:N	2.11	0.65
1:C:384:VAL:HG22	1:C:410:MET:HB3	1.83	0.60
1:B:46:ILE:HD12	1:B:344:SER:HB3	1.86	0.58
1:A:285:ASP:OD2	1:C:55:PRO:O	2.22	0.57
1:C:172:TRP:HB2	1:C:189:THR:HB	1.85	0.56
1:B:70:GLU:HG2	1:B:417:LYS:HD3	1.88	0.56
1:B:447:GLU:HA	1:B:448:ASP:C	2.27	0.55
1:C:380:GLY:HA2	5:C:898:HOH:O	2.07	0.54
1:B:475:GLU:HB2	1:B:478:GLN:HB2	1.89	0.54
1:D:46:ILE:HD12	1:D:344:SER:HB3	1.89	0.54
1:C:46:ILE:HD12	1:C:344:SER:HB3	1.90	0.53
1:C:447:GLU:HG2	1:C:448:ASP:HA	1.90	0.53
1:C:323:LEU:HG	1:C:351:ILE:HD12	1.91	0.52
1:D:55:PRO:O	1:D:58:GLN:HG2	2.10	0.51
1:A:172:TRP:HB2	1:A:189:THR:HB	1.92	0.51
1:D:172:TRP:HB2	1:D:189:THR:HB	1.93	0.51
1:B:388:GLY:HA3	1:B:401:LEU:HD12	1.92	0.50
1:B:323:LEU:HG	1:B:351:ILE:HD12	1.94	0.49
1:A:388:GLY:HA3	1:A:401:LEU:HD12	1.95	0.49
1:D:121:ASP:OD1	1:D:122:PRO:HD2	2.11	0.49
1:B:239:GLN:HA	1:B:242:ILE:HD12	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:362:LYS:HE2	5:D:560:HOH:O	2.13	0.48
1:C:112:TYR:CZ	1:C:136:LYS:HE3	2.49	0.48
1:D:323:LEU:HG	1:D:351:ILE:HD12	1.95	0.47
1:D:54:ILE:N	1:D:55:PRO:HD2	2.30	0.47
1:B:433:ASN:HB3	1:B:458:LEU:HD22	1.97	0.47
1:C:82:LYS:HE2	1:C:88:THR:HB	1.96	0.47
1:C:239:GLN:HA	1:C:242:ILE:HD12	1.97	0.47
1:A:70:GLU:HG2	1:A:417:LYS:HD3	1.97	0.46
1:A:157:LYS:HD3	1:A:168:GLN:HB3	1.97	0.46
1:B:75:ASN:HD21	1:B:449:ASN:H	1.65	0.45
1:C:323:LEU:HD11	1:C:367:THR:HB	1.98	0.45
1:A:46:ILE:HD12	1:A:344:SER:HB3	1.98	0.45
1:A:455:PRO:HA	1:A:476:GLN:HA	1.99	0.45
1:B:170:GLN:HB2	1:B:191:TYR:HD2	1.82	0.45
1:D:142:ILE:HA	1:D:145:TRP:HD1	1.82	0.44
1:A:200:THR:HG23	1:A:229:ASP:HB3	2.00	0.44
1:A:119:ALA:HB3	1:A:131:TYR:CD2	2.52	0.44
1:D:297:ASN:N	1:D:298:PRO:HD2	2.32	0.44
1:B:449:ASN:OD1	1:B:449:ASN:N	2.50	0.44
1:C:326:ILE:HG22	1:C:337:VAL:HA	1.99	0.44
1:D:388:GLY:HA3	1:D:401:LEU:HD12	2.00	0.44
1:D:447:GLU:HG3	1:D:448:ASP:OD1	2.18	0.43
1:C:55:PRO:O	1:C:58:GLN:HG2	2.19	0.43
1:C:238:VAL:O	1:C:242:ILE:HG13	2.19	0.42
1:B:119:ALA:HB3	1:B:131:TYR:CD2	2.55	0.42
1:B:70:GLU:CG	1:B:417:LYS:HD3	2.48	0.42
1:D:447:GLU:HG3	1:D:448:ASP:HA	2.01	0.42
1:B:157:LYS:HD3	1:B:168:GLN:HB3	2.00	0.42
1:B:172:TRP:HB2	1:B:189:THR:HB	2.01	0.42
1:B:355:ASN:HB3	1:B:366:PHE:HB2	2.01	0.42
1:D:54:ILE:N	1:D:55:PRO:CD	2.83	0.42
1:C:203:THR:HB	1:C:227:ILE:HD11	2.01	0.42
1:B:97:TRP:CD1	1:B:175:SER:HA	2.55	0.41
1:B:410:MET:HE1	1:B:419:PHE:CE1	2.55	0.41
1:D:389:TYR:CD1	1:D:400:PRO:HA	2.56	0.41
1:C:270:VAL:HG12	1:C:354:ALA:HB1	2.02	0.41
1:C:399:LYS:HD2	1:C:399:LYS:HA	1.86	0.41
1:A:323:LEU:HG	1:A:351:ILE:HD12	2.03	0.41
1:B:190:ASP:HB3	1:B:200:THR:HB	2.02	0.41
1:A:475:GLU:HB3	5:A:651:HOH:O	2.20	0.40
1:C:388:GLY:HA3	1:C:401:LEU:HD12	2.04	0.40

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	446/456 (98%)	428 (96%)	18 (4%)	0	100	100
1	B	448/456 (98%)	430 (96%)	18 (4%)	0	100	100
1	C	446/456 (98%)	429 (96%)	17 (4%)	0	100	100
1	D	450/456 (99%)	434 (96%)	15 (3%)	1 (0%)	52	43
All	All	1790/1824 (98%)	1721 (96%)	68 (4%)	1 (0%)	56	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	125	SER

### 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	389/395 (98%)	383 (98%)	6 (2%)	72	68
1	B	391/395 (99%)	385 (98%)	6 (2%)	72	68
1	C	389/395 (98%)	384 (99%)	5 (1%)	76	72
1	D	393/395 (100%)	382 (97%)	11 (3%)	51	39
All	All	1562/1580 (99%)	1534 (98%)	28 (2%)	66	60

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	ASP
1	A	166	LYS
1	A	214	ASN
1	A	255	LEU
1	A	367	THR
1	A	443	ARG
1	B	56	GLN
1	B	82	LYS
1	B	95	ASP
1	B	140	LYS
1	B	248	ASP
1	B	255	LEU
1	C	82	LYS
1	C	95	ASP
1	C	255	LEU
1	C	314	GLU
1	C	481	VAL
1	D	39	ASN
1	D	82	LYS
1	D	95	ASP
1	D	124	ASP
1	D	166	LYS
1	D	233	LYS
1	D	244	GLU
1	D	248	ASP
1	D	255	LEU
1	D	447	GLU
1	D	481	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	50	ASN
1	C	402	ASN
1	D	39	ASN
1	D	126	ASN

### 5.3.3 RNA ⓘ

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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	SO4	A	485	-	4,4,4	0.59	0	6,6,6	0.37	0
3	SO4	B	3	-	4,4,4	0.49	0	6,6,6	0.42	0
3	SO4	B	4	-	4,4,4	0.32	0	6,6,6	0.27	0
3	SO4	B	5	-	4,4,4	0.51	0	6,6,6	0.21	0
4	PGE	C	485	-	9,9,9	0.34	0	8,8,8	0.50	0
3	SO4	D	2	-	4,4,4	0.61	0	6,6,6	0.33	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	SO4	A	485	-	-	0/0/0/0	0/0/0/0
3	SO4	B	3	-	-	0/0/0/0	0/0/0/0
3	SO4	B	4	-	-	0/0/0/0	0/0/0/0
3	SO4	B	5	-	-	0/0/0/0	0/0/0/0
4	PGE	C	485	-	-	0/7/7/7	0/0/0/0
3	SO4	D	2	-	-	0/0/0/0	0/0/0/0

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

### 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, 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	448/456 (98%)	0.93	31 (6%)	20	30	7, 13, 21, 26	5 (1%)
1	B	448/456 (98%)	0.96	45 (10%)	9	15	3, 13, 23, 26	5 (1%)
1	C	448/456 (98%)	0.86	22 (4%)	33	45	6, 13, 19, 24	5 (1%)
1	D	448/456 (98%)	0.95	41 (9%)	11	18	4, 13, 21, 26	5 (1%)
All	All	1792/1824 (98%)	0.93	139 (7%)	16	25	3, 13, 21, 26	20 (1%)

All (139) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	125	SER	8.1
1	D	84	ALA	7.7
1	D	81	GLY	7.2
1	B	84	ALA	7.0
1	A	446	TYR	6.2
1	B	125	SER	6.1
1	D	122	PRO	6.0
1	B	446	TYR	5.5
1	B	448	ASP	5.3
1	D	88	THR	5.2
1	D	85	SER	5.0
1	A	444	GLY	5.0
1	C	123	LYS	5.0
1	C	84	ALA	4.9
1	D	82	LYS	4.9
1	B	128	THR	4.8
1	A	447	GLU	4.7
1	C	122	PRO	4.7
1	D	125	SER	4.6
1	B	124	ASP	4.5
1	B	86	GLY	4.5

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Mol	Chain	Res	Type	RSRZ
1	C	125	SER	4.4
1	D	83	ASN	4.4
1	D	87	ASN	4.4
1	A	445	PHE	4.3
1	C	247	ALA	4.3
1	A	196	TYR	4.2
1	D	196	TYR	4.1
1	D	138	GLY	4.1
1	B	85	SER	4.0
1	A	443	ARG	4.0
1	C	467	SER	4.0
1	B	310	GLY	4.0
1	B	445	PHE	3.9
1	B	88	THR	3.9
1	B	126	ASN	3.9
1	D	123	LYS	3.8
1	B	309[A]	GLU	3.8
1	D	86	GLY	3.7
1	B	87	ASN	3.7
1	D	246	GLY	3.6
1	B	447	GLU	3.6
1	B	249	THR	3.5
1	B	83	ASN	3.5
1	D	128	THR	3.5
1	A	122	PRO	3.4
1	B	127	ASP	3.3
1	D	142	ILE	3.3
1	A	211	PRO	3.3
1	D	34	SER	3.2
1	C	87	ASN	3.2
1	A	448	ASP	3.2
1	B	158	PHE	3.2
1	C	85	SER	3.2
1	D	120	GLY	3.1
1	D	300	PHE	3.1
1	B	467	SER	3.1
1	D	249	THR	3.0
1	C	124	ASP	3.0
1	D	245	GLY	3.0
1	B	247	ALA	3.0
1	B	363	TRP	3.0
1	B	242	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	140	LYS	2.9
1	A	121	ASP	2.9
1	D	307	LEU	2.9
1	D	247	ALA	2.9
1	B	123	LYS	2.9
1	A	128	THR	2.9
1	A	467	SER	2.8
1	B	153	LYS	2.8
1	D	467	SER	2.8
1	C	86	GLY	2.8
1	B	89	ILE	2.8
1	B	311	SER	2.8
1	B	121	ASP	2.7
1	B	245	GLY	2.7
1	D	126	ASN	2.7
1	C	81	GLY	2.6
1	C	83	ASN	2.6
1	B	133	PHE	2.6
1	A	123	LYS	2.6
1	A	357	PHE	2.5
1	D	473	VAL	2.5
1	B	186	LEU	2.5
1	D	308	LEU	2.5
1	D	143	ASP	2.5
1	B	122	PRO	2.5
1	A	132	LEU	2.4
1	C	89	ILE	2.4
1	A	247	ALA	2.4
1	D	145	TRP	2.4
1	A	108	THR	2.4
1	D	337	VAL	2.4
1	B	154	ASP	2.4
1	B	264	ASN	2.4
1	B	82	LYS	2.4
1	D	181	ASP	2.4
1	B	206	VAL	2.4
1	A	442	ASN	2.3
1	B	208	MET	2.3
1	B	166	LYS	2.3
1	A	124	ASP	2.3
1	C	268	TYR	2.3
1	B	252	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	255	LEU	2.3
1	B	307	LEU	2.3
1	D	139	ASP	2.3
1	D	474	LEU	2.3
1	D	127	ASP	2.3
1	A	165	LEU	2.2
1	D	314[A]	GLU	2.2
1	A	111	GLY	2.2
1	B	246	GLY	2.2
1	C	213	ASP	2.2
1	D	461	ILE	2.2
1	A	158	PHE	2.2
1	D	124	ASP	2.1
1	D	472	ARG	2.1
1	B	117	ALA	2.1
1	C	354	ALA	2.1
1	A	172	TRP	2.1
1	D	165	LEU	2.1
1	A	161	ASN	2.1
1	B	308	LEU	2.1
1	C	387	LEU	2.1
1	A	184	VAL	2.1
1	D	195	GLN	2.1
1	B	443	ARG	2.1
1	C	445	PHE	2.1
1	B	181	ASP	2.1
1	C	148	ALA	2.1
1	A	80	LYS	2.1
1	C	194	LYS	2.1
1	C	196	TYR	2.0
1	C	193	GLY	2.0
1	A	126	ASN	2.0
1	A	144	SER	2.0
1	A	461	ILE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	SO4	B	5	5/5	0.94	0.21	2.96	31,31,31,31	0
3	SO4	B	4	5/5	0.82	0.21	1.35	33,33,33,33	0
4	PGE	C	485	10/10	0.87	0.15	0.05	20,20,21,21	0
3	SO4	D	2	5/5	0.93	0.12	-0.66	31,31,32,32	0
3	SO4	A	485	5/5	0.95	0.10	-1.42	30,30,30,30	0
2	CA	C	1	1/1	0.97	0.08	-2.39	10,10,10,10	0
2	CA	B	1	1/1	0.99	0.04	-2.63	12,12,12,12	0
2	CA	A	1	1/1	0.99	0.05	-2.67	13,13,13,13	0
2	CA	D	1	1/1	0.99	0.07	-4.20	14,14,14,14	0
3	SO4	B	3	5/5	0.93	0.15	-	32,32,32,32	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.