



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2016 – 04:13 PM EDT

PDB ID : 5G2P  
Title : The crystal structure of a S-selective transaminase from *Arthrobacter* sp.  
Authors : van Oosterwijk, N.; Willies, S.; Hekelaar, J.; Terwisscha van Scheltinga, A.C.;  
Turner, N.J.; Dijkstra, B.W.  
Deposited on : 2016-04-12  
Resolution : 1.89 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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027939  
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) : rb-20027939

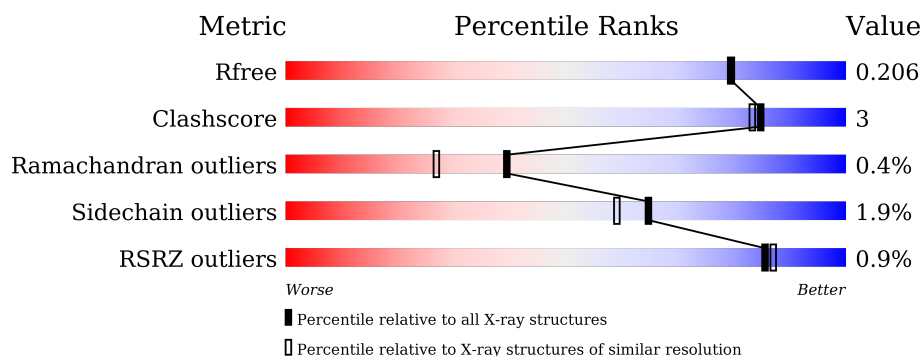
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.89 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	485	<div> <div>87%</div> <div>6% • 5%</div> </div>
1	B	485	<div> <div>%</div> <div>89%</div> <div>6% •</div> </div>
1	C	485	<div> <div>%</div> <div>86%</div> <div>8% • 5%</div> </div>
1	D	485	<div> <div>%</div> <div>88%</div> <div>8% 5%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15740 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.

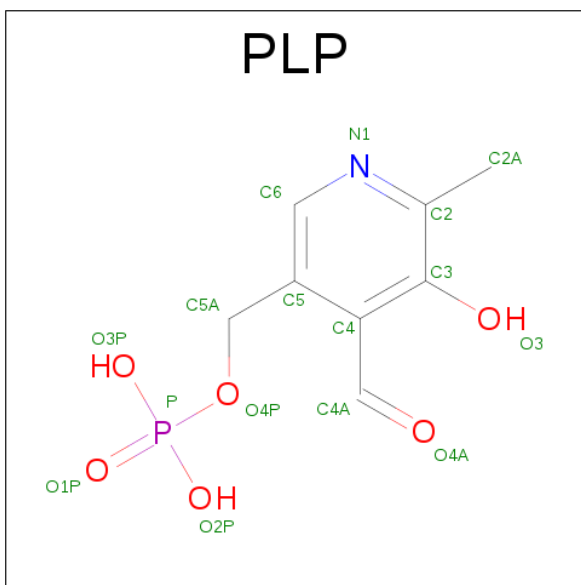
- Molecule 1 is a protein called TRANSAMINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	461	Total	C	N	O	S	0	0	0
			3624	2297	617	692	18			
1	B	464	Total	C	N	O	S	0	0	0
			3641	2305	620	698	18			
1	C	461	Total	C	N	O	S	0	0	0
			3619	2294	616	691	18			
1	D	462	Total	C	N	O	S	0	0	0
			3627	2298	617	694	18			

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C<sub>8</sub>H<sub>10</sub>NO<sub>6</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	C	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	D	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

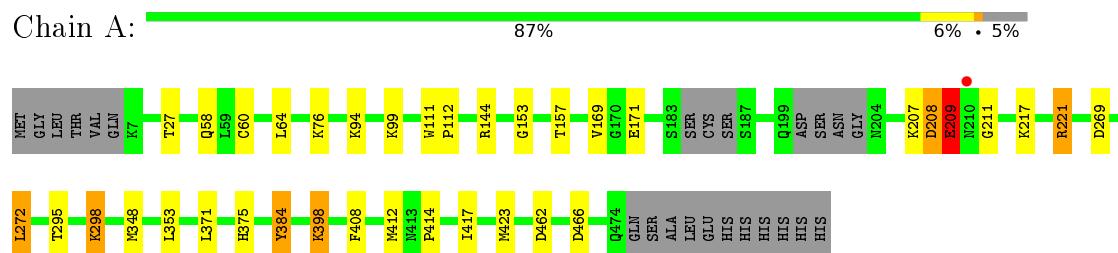
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	360	Total	O	0	0
			360	360		
4	B	318	Total	O	0	0
			318	318		
4	C	269	Total	O	0	0
			269	269		
4	D	218	Total	O	0	0
			218	218		

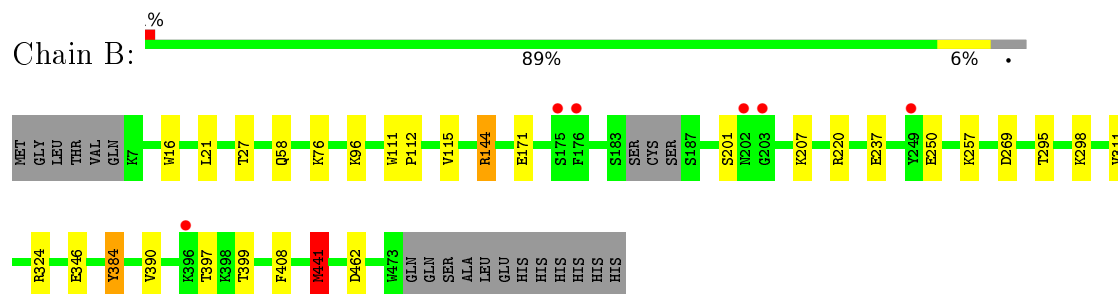
### 3 Residue-property plots [i](#)

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.

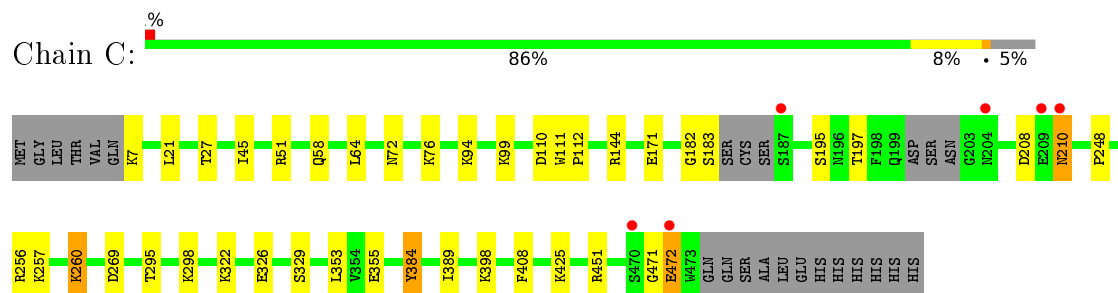
#### • Molecule 1: TRANSAMINASE



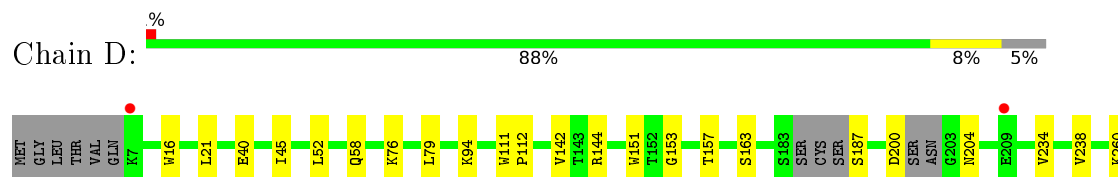
#### • Molecule 1: TRANSAMINASE



#### • Molecule 1: TRANSAMINASE



#### • Molecule 1: TRANSAMINASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	64.40 Å 142.66 Å 99.68 Å 90.00° 105.33° 90.00°	Depositor
Resolution (Å)	46.84 – 1.89 46.84 – 1.89	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.84-1.89) 99.8 (46.84-1.89)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.92 (at 1.90 Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, $R_{free}$	0.157 , 0.200 0.161 , 0.206	Depositor DCC
$R_{free}$ test set	6859 reflections (5.27%)	DCC
Wilson B-factor (Å <sup>2</sup> )	18.2	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 46.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	15740	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, PLP

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	0.59	0/3700	0.68	1/5009 (0.0%)
1	B	0.59	0/3718	0.69	2/5035 (0.0%)
1	C	0.54	0/3695	0.65	0/5002
1	D	0.54	0/3703	0.67	0/5013
All	All	0.56	0/14816	0.67	3/20059 (0.0%)

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	A	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	348	MET	CG-SD-CE	-5.84	90.86	100.20
1	B	441	MET	CG-SD-CE	-5.82	90.89	100.20
1	B	462	ASP	CB-CG-OD1	5.04	122.84	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	209	GLU	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	3624	0	3571	32	0
1	B	3641	0	3582	16	0
1	C	3619	0	3566	26	0
1	D	3627	0	3570	23	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	15	0	7	1	0
3	B	15	0	7	1	0
3	C	15	0	6	1	0
3	D	15	0	6	1	0
4	A	360	0	0	3	0
4	B	318	0	0	1	0
4	C	269	0	0	1	0
4	D	218	0	0	1	0
All	All	15740	0	14315	85	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 3.

All (85) 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:A:221:ARG:HG3	1:A:221:ARG:HH11	1.05	1.09
1:A:221:ARG:HG3	1:A:221:ARG:NH1	1.85	0.86
1:A:423:MET:CE	4:A:3339:HOH:O	2.21	0.86
1:B:397:THR:OG1	1:B:399:THR:HG22	1.80	0.81
1:C:322:LYS:HE3	4:C:3206:HOH:O	1.82	0.79
1:C:471:GLY:O	1:C:472:GLU:CB	2.32	0.77
1:B:220:ARG:HD3	4:B:3186:HOH:O	1.85	0.76
1:B:390:VAL:HB	1:B:441:MET:HG3	1.70	0.73
1:C:471:GLY:O	1:C:472:GLU:HB2	1.92	0.70
1:A:423:MET:HE3	4:A:3339:HOH:O	1.91	0.67
1:B:207:LYS:NZ	1:B:250:GLU:OE2	2.26	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:MET:HE2	4:A:3339:HOH:O	1.88	0.66
1:A:221:ARG:HH11	1:A:221:ARG:CG	1.94	0.64
1:C:27:THR:HG21	1:C:171:GLU:HB2	1.79	0.64
1:C:76:LYS:HG2	1:D:76:LYS:HG2	1.86	0.58
1:A:371:LEU:HD21	1:A:462:ASP:HB2	1.86	0.57
1:D:399:THR:HG22	1:D:400:PRO:HD2	1.87	0.56
1:C:182:GLY:O	1:C:183:SER:CB	2.54	0.56
1:A:76:LYS:HG2	1:B:76:LYS:HG2	1.87	0.56
1:C:21:LEU:HD13	1:D:94:LYS:HE3	1.89	0.54
1:D:200:ASP:HB2	1:D:204:ASN:H	1.74	0.53
1:A:94:LYS:HE3	1:B:21:LEU:HD13	1.92	0.52
1:D:153:GLY:O	1:D:157:THR:HG23	2.11	0.51
1:A:412:MET:HG3	1:A:417:ILE:HD11	1.92	0.51
1:B:298:LYS:NZ	3:B:2000:PLP:O3	2.43	0.50
1:B:115:VAL:HG22	1:B:311:VAL:HG22	1.94	0.49
1:A:221:ARG:HH12	1:C:197:THR:HG23	1.77	0.49
1:D:298:LYS:NZ	3:D:2000:PLP:O3	2.46	0.49
1:A:209:GLU:C	1:A:211:GLY:H	2.15	0.49
1:B:269:ASP:HA	1:B:295:THR:OG1	2.12	0.49
1:A:209:GLU:C	1:A:211:GLY:N	2.65	0.48
1:A:153:GLY:O	1:A:157:THR:HG23	2.13	0.48
1:A:60:CYS:CB	1:A:272:LEU:HD21	2.43	0.48
1:A:269:ASP:HA	1:A:295:THR:OG1	2.13	0.48
1:C:269:ASP:HA	1:C:295:THR:OG1	2.13	0.47
1:C:45:ILE:HD12	1:C:51:ARG:HG2	1.96	0.47
1:D:260:LYS:HE3	4:D:3068:HOH:O	2.15	0.47
1:C:329:SER:HB2	1:D:151:TRP:NE1	2.30	0.47
1:C:471:GLY:O	1:C:472:GLU:HB3	2.15	0.46
1:D:111:TRP:N	1:D:112:PRO:CD	2.78	0.46
1:A:111:TRP:N	1:A:112:PRO:CD	2.79	0.46
1:C:256:ARG:CZ	1:C:260:LYS:HE2	2.47	0.45
1:C:195:SER:HB2	1:C:248:PRO:HD2	1.98	0.45
1:D:329:SER:HB3	1:D:332:ALA:HB2	1.98	0.45
1:A:60:CYS:HB2	1:A:272:LEU:HD21	1.97	0.45
1:D:366:SER:O	1:D:370:LEU:HG	2.16	0.45
1:C:182:GLY:O	1:C:183:SER:HB2	2.16	0.45
1:C:208:ASP:OD1	1:C:210:ASN:HB2	2.16	0.44
1:A:169:VAL:HG23	1:B:324:ARG:HB2	1.99	0.44
1:C:99:LYS:HG3	1:D:16:TRP:CE2	2.53	0.44
1:A:375:HIS:HE1	1:A:466:ASP:OD1	2.00	0.44
1:D:40:GLU:HG2	1:D:45:ILE:HD11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:THR:HG21	1:A:171:GLU:HB2	2.00	0.43
1:D:200:ASP:HB2	1:D:204:ASN:N	2.32	0.43
1:A:209:GLU:CA	1:A:211:GLY:H	2.31	0.43
1:C:298:LYS:NZ	3:C:2000:PLP:O3	2.50	0.43
1:D:142:VAL:HB	1:D:234:VAL:HG22	2.00	0.43
1:C:425:LYS:HD2	1:C:425:LYS:HA	1.73	0.43
1:C:94:LYS:HE3	1:D:21:LEU:HD13	2.01	0.43
1:A:208:ASP:OD2	1:A:217:LYS:HG3	2.19	0.43
1:A:221:ARG:CG	1:A:221:ARG:NH1	2.62	0.42
1:A:414:PRO:HA	1:A:417:ILE:HD12	2.01	0.42
1:D:111:TRP:N	1:D:112:PRO:HD3	2.35	0.42
1:C:111:TRP:N	1:C:112:PRO:CD	2.83	0.42
1:D:271:VAL:O	1:D:298:LYS:HB2	2.20	0.42
1:A:64:LEU:HD12	1:A:353:LEU:HD12	2.01	0.42
1:C:384:TYR:CD1	1:C:384:TYR:N	2.89	0.41
1:D:238:VAL:HG21	1:D:283:GLY:HA3	2.02	0.41
1:A:111:TRP:N	1:A:112:PRO:HD3	2.35	0.41
1:A:398:LYS:HD2	1:A:398:LYS:HA	1.76	0.41
1:C:64:LEU:HD12	1:C:353:LEU:HD12	2.03	0.41
1:A:207:LYS:HE3	1:A:211:GLY:O	2.20	0.41
1:B:27:THR:HG21	1:B:171:GLU:HB2	2.01	0.41
1:C:72:ASN:ND2	1:D:79:LEU:O	2.53	0.41
1:B:384:TYR:N	1:B:384:TYR:CD1	2.87	0.41
1:A:99:LYS:HG3	1:B:16:TRP:CE2	2.55	0.41
1:D:413:ASN:HA	1:D:414:PRO:HD2	1.93	0.41
1:A:298:LYS:NZ	3:A:2000:PLP:O3	2.55	0.40
1:B:96:LYS:HE2	1:B:346:GLU:OE2	2.21	0.40
1:C:99:LYS:C	1:C:99:LYS:HD3	2.42	0.40
1:D:414:PRO:HA	1:D:417:ILE:HD12	2.04	0.40
1:A:384:TYR:CD1	1:A:384:TYR:N	2.89	0.40
1:C:326:GLU:HB3	1:D:163:SER:O	2.22	0.40
1:B:111:TRP:N	1:B:112:PRO:HD3	2.36	0.40
1:B:144:ARG:NH2	1:B:237:GLU:H	2.20	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	455/485 (94%)	430 (94%)	23 (5%)	2 (0%)	39	27
1	B	460/485 (95%)	439 (95%)	20 (4%)	1 (0%)	52	42
1	C	455/485 (94%)	432 (95%)	20 (4%)	3 (1%)	26	14
1	D	456/485 (94%)	437 (96%)	18 (4%)	1 (0%)	52	42
All	All	1826/1940 (94%)	1738 (95%)	81 (4%)	7 (0%)	39	27

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	210	ASN
1	C	472	GLU
1	A	298	LYS
1	B	58	GLN
1	C	58	GLN
1	A	58	GLN
1	D	58	GLN

### 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	388/409 (95%)	380 (98%)	8 (2%)	61	55
1	B	390/409 (95%)	384 (98%)	6 (2%)	72	69
1	C	387/409 (95%)	376 (97%)	11 (3%)	51	41

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	388/409 (95%)	383 (99%)	5 (1%)	76	73
All	All	1553/1636 (95%)	1523 (98%)	30 (2%)	65	59

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	144	ARG
1	A	208	ASP
1	A	209	GLU
1	A	221	ARG
1	A	272	LEU
1	A	384	TYR
1	A	398	LYS
1	A	408	PHE
1	B	144	ARG
1	B	201	SER
1	B	257	LYS
1	B	384	TYR
1	B	408	PHE
1	B	441	MET
1	C	7	LYS
1	C	110	ASP
1	C	144	ARG
1	C	257	LYS
1	C	260	LYS
1	C	355	GLU
1	C	384	TYR
1	C	389	ILE
1	C	398	LYS
1	C	408	PHE
1	C	451	ARG
1	D	52	LEU
1	D	144	ARG
1	D	187	SER
1	D	384	TYR
1	D	408	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	372	GLN

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Mol	Chain	Res	Type
1	B	290	GLN
1	B	359	ASN
1	C	290	GLN
1	C	416	GLN

### 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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	PLP	A	2000	1	15,15,16	3.73	3 (20%)	21,22,23	2.18	6 (28%)
3	PLP	B	2000	1	15,15,16	3.57	3 (20%)	21,22,23	2.47	4 (19%)
3	PLP	C	2000	1	15,15,16	3.39	3 (20%)	21,22,23	2.37	5 (23%)
3	PLP	D	2000	1	15,15,16	4.00	3 (20%)	21,22,23	2.50	5 (23%)

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	PLP	A	2000	1	-	0/6/6/8	0/1/1/1
3	PLP	B	2000	1	-	0/6/6/8	0/1/1/1
3	PLP	C	2000	1	-	0/6/6/8	0/1/1/1
3	PLP	D	2000	1	-	0/6/6/8	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	2000	PLP	C3-C4	3.10	1.47	1.40
3	D	2000	PLP	C3-C4	3.45	1.47	1.40
3	B	2000	PLP	C3-C4	3.56	1.48	1.40
3	A	2000	PLP	C3-C4	3.90	1.48	1.40
3	C	2000	PLP	C5-C4	8.15	1.50	1.40
3	B	2000	PLP	C5-C4	8.66	1.50	1.40
3	A	2000	PLP	C5-C4	8.67	1.50	1.40
3	D	2000	PLP	C5-C4	8.99	1.51	1.40
3	C	2000	PLP	C3-C2	9.59	1.47	1.40
3	B	2000	PLP	C3-C2	9.79	1.47	1.40
3	A	2000	PLP	C3-C2	10.43	1.48	1.40
3	D	2000	PLP	C3-C2	11.60	1.49	1.40

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	2000	PLP	C3-C4-C5	-3.96	114.04	118.68
3	B	2000	PLP	C4A-C4-C3	-3.87	113.68	120.42
3	C	2000	PLP	C3-C4-C5	-3.50	114.58	118.68
3	A	2000	PLP	C4A-C4-C3	-2.80	115.55	120.42
3	C	2000	PLP	C4A-C4-C3	-2.44	116.17	120.42
3	D	2000	PLP	C4A-C4-C3	-2.40	116.24	120.42
3	A	2000	PLP	C3-C4-C5	-2.39	115.87	118.68
3	A	2000	PLP	C2A-C2-C3	-2.36	118.51	120.90
3	A	2000	PLP	O4P-P-O1P	-2.16	101.65	107.08
3	A	2000	PLP	O3-C3-C2	2.07	120.57	117.53
3	C	2000	PLP	O3P-P-O1P	2.42	118.51	110.63
3	B	2000	PLP	C6-N1-C2	2.55	124.38	119.26
3	D	2000	PLP	C6-C5-C4	2.97	120.64	118.12
3	C	2000	PLP	O3-C3-C2	3.07	122.03	117.53
3	D	2000	PLP	O3-C3-C2	3.41	122.53	117.53
3	B	2000	PLP	O3-C3-C2	3.85	123.17	117.53
3	A	2000	PLP	C4A-C4-C5	7.38	128.50	120.90
3	C	2000	PLP	C4A-C4-C5	8.06	129.19	120.90
3	B	2000	PLP	C4A-C4-C5	8.28	129.42	120.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	2000	PLP	C4A-C4-C5	8.50	129.65	120.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2000	PLP	1	0
3	B	2000	PLP	1	0
3	C	2000	PLP	1	0
3	D	2000	PLP	1	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 ⓘ

### 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	461/485 (95%)	-0.49	1 (0%) 95 95	12, 19, 36, 72	0
1	B	464/485 (95%)	-0.45	6 (1%) 79 82	12, 18, 34, 63	0
1	C	461/485 (95%)	-0.30	6 (1%) 79 82	15, 23, 42, 75	0
1	D	462/485 (95%)	-0.31	4 (0%) 85 87	15, 23, 39, 59	0
All	All	1848/1940 (95%)	-0.39	17 (0%) 85 87	12, 21, 39, 75	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	176	PHE	4.9
1	C	209	GLU	4.8
1	C	210	ASN	4.7
1	B	202	ASN	3.0
1	D	473	TRP	2.9
1	B	175	SER	2.9
1	B	249	TYR	2.9
1	D	209	GLU	2.8
1	C	187	SER	2.5
1	A	210	ASN	2.4
1	B	396	LYS	2.4
1	C	472	GLU	2.3
1	C	204	ASN	2.2
1	D	472	GLU	2.2
1	B	203	GLY	2.1
1	C	470	SER	2.1
1	D	7	LYS	2.1

### 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
2	CL	B	1474	1/1	0.99	0.08	0.69	23,23,23,23	0
2	CL	A	1475	1/1	0.99	0.07	0.57	25,25,25,25	0
3	PLP	B	2000	15/16	0.98	0.11	0.28	12,14,17,19	0
3	PLP	C	2000	15/16	0.98	0.12	0.16	15,19,24,25	0
2	CL	D	1474	1/1	0.99	0.08	-0.03	25,25,25,25	0
3	PLP	D	2000	15/16	0.98	0.12	-0.07	17,18,21,21	0
2	CL	C	1474	1/1	0.99	0.07	-0.16	26,26,26,26	0
3	PLP	A	2000	15/16	0.99	0.10	-0.33	15,16,18,20	0

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

There are no such residues in this entry.