



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:31 AM GMT

PDB ID : 3EY9  
Title : Structural basis for membrane binding and catalytic activation of the peripheral membrane enzyme pyruvate oxidase from Escherichia coli  
Authors : Neumann, P.; Weidner, A.; Pech, A.; Stubbs, M.T.; Tittmann, K.  
Deposited on : 2008-10-20  
Resolution : 2.90 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

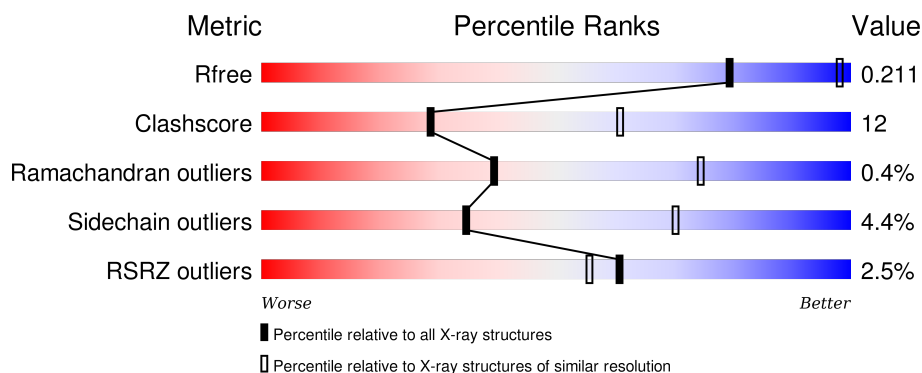
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	572	<div> <div>3%</div> <div>77%</div> <div>22%</div> <div>.</div> </div>
1	B	572	<div> <div>2%</div> <div>76%</div> <div>22%</div> <div>.</div> </div>

## 2 Entry composition [i](#)

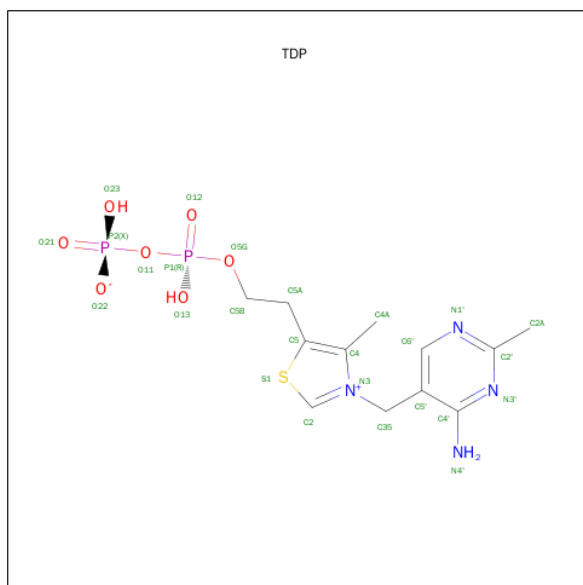
There are 5 unique types of molecules in this entry. The entry contains 8854 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Pyruvate dehydrogenase [cytochrome].

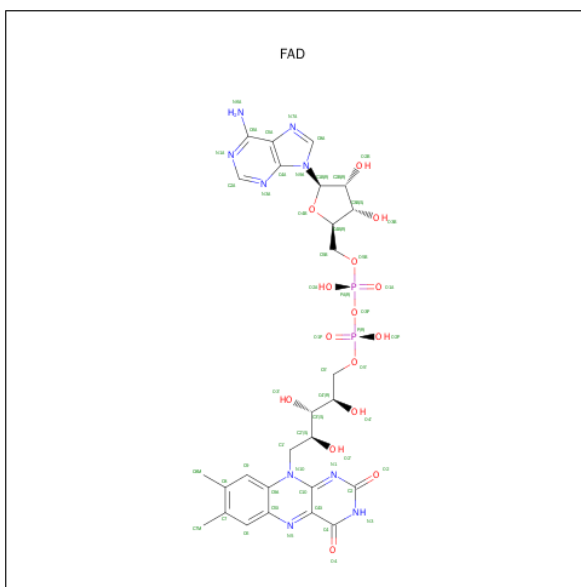
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	571	Total	C	N	O	S	0	0	0
			4342	2750	753	810	29			
1	B	571	Total	C	N	O	S	0	0	0
			4342	2750	753	810	29			

- Molecule 2 is THIAMIN DIPHOSPHATE (three-letter code: TDP) (formula:  $C_{12}H_{18}N_4O_7P_2S$ ).



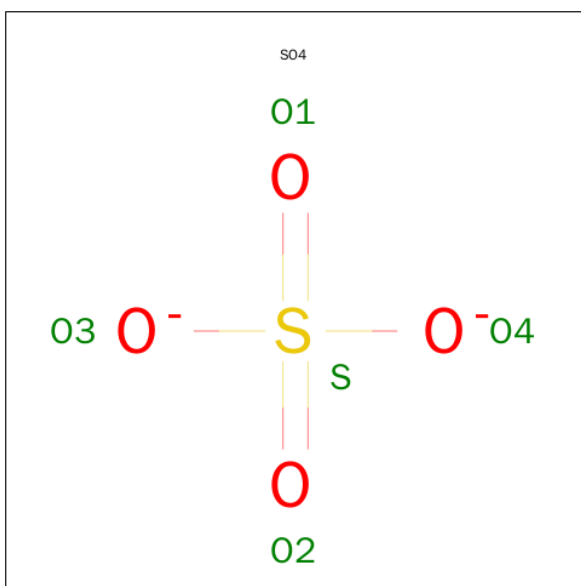
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P S	0	0
			26	12	4	7	2 1		
2	B	1	Total	C	N	O	P S	0	0
			26	12	4	7	2 1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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



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

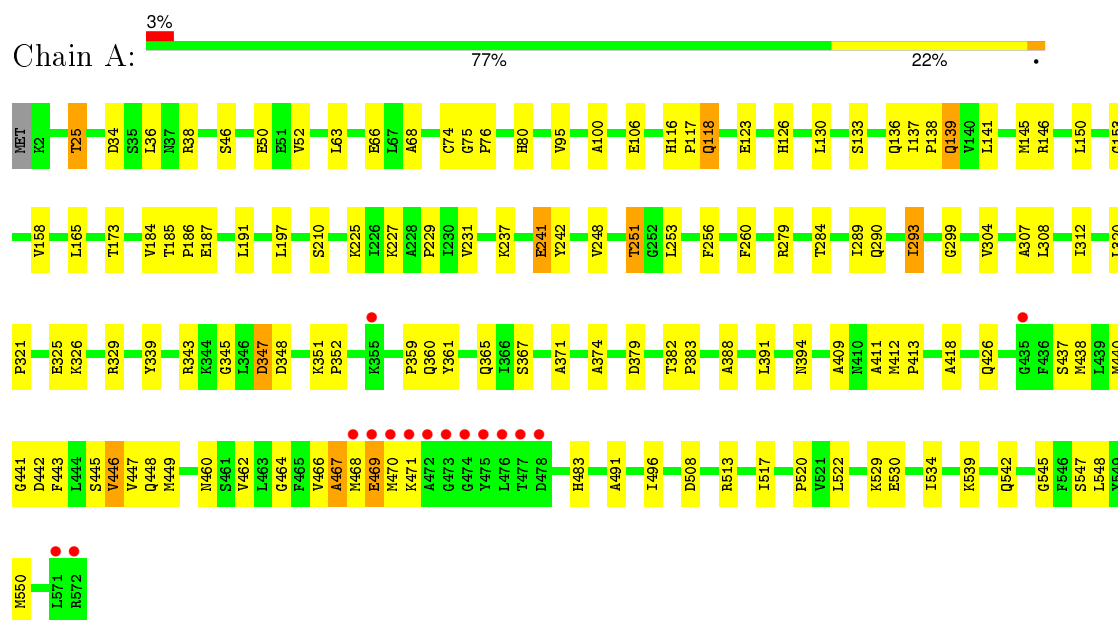
- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mg	0	0
			1	1		
5	A	1	Total	Mg	0	0
			1	1		

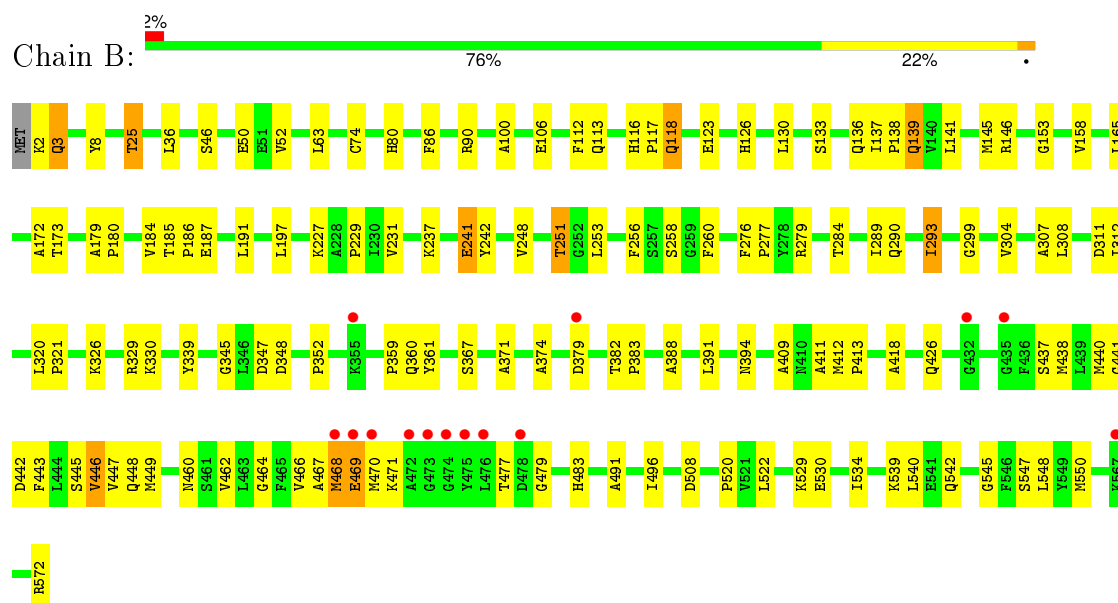
### 3 Residue-property plots

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: Pyruvate dehydrogenase [cytochrome]



#### • Molecule 1: Pyruvate dehydrogenase [cytochrome]



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	151.37Å 151.37Å 153.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.97 – 2.90 29.97 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.97-2.90) 98.8 (29.97-2.90)	Depositor EDS
$R_{merge}$	0.01	Depositor
$R_{sym}$	0.01	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.75 (at 2.90Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, $R_{free}$	0.182 , 0.216 0.180 , 0.211	Depositor DCC
$R_{free}$ test set	1982 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	73.5	Xtriage
Anisotropy	0.427	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 40.7	EDS
Estimated twinning fraction	0.010 for -h,l,k 0.000 for -l,-k,-h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 39657 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8854	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, TDP, SO4, FAD

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	0.37	0/4430	0.54	0/6001
1	B	0.36	0/4430	0.54	0/6001
All	All	0.37	0/8860	0.54	0/12002

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

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	4342	0	4348	101	0
1	B	4342	0	4348	110	0
2	A	26	0	16	6	0
2	B	26	0	16	7	0
3	A	53	0	31	2	0
3	B	53	0	31	3	0
4	A	5	0	0	1	0
4	B	5	0	0	1	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
All	All	8854	0	8790	203	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 12.

All (203) 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)
2:A:611:TDP:C2	2:A:611:TDP:H2	0.97	1.49
2:B:611:TDP:H2	2:B:611:TDP:C2	0.97	1.49
1:A:466:VAL:HG21	1:A:529:LYS:HB3	1.45	0.98
1:B:466:VAL:HG21	1:B:529:LYS:HB3	1.45	0.94
1:B:547:SER:H	1:B:550:MET:HE3	1.35	0.91
1:A:371:ALA:H	1:A:426:GLN:NE2	1.69	0.89
1:A:371:ALA:H	1:A:426:GLN:HE22	1.19	0.87
1:A:123:GLU:OE2	1:B:116:HIS:HD2	1.57	0.86
1:B:371:ALA:H	1:B:426:GLN:NE2	1.73	0.86
1:B:371:ALA:H	1:B:426:GLN:HE22	1.27	0.82
1:B:330:LYS:HZ3	1:B:572:ARG:HD3	1.45	0.81
1:A:116:HIS:HD2	1:B:123:GLU:OE2	1.63	0.80
1:A:539:LYS:HA	1:A:542:GLN:HE21	1.48	0.79
1:A:227:LYS:HG2	1:A:329:ARG:HB3	1.64	0.79
1:B:118:GLN:HG2	1:B:130:LEU:HB2	1.66	0.77
1:A:445:SER:HB3	1:A:449:MET:HE3	1.66	0.77
1:B:445:SER:HB3	1:B:449:MET:HE3	1.66	0.77
1:B:227:LYS:HG2	1:B:329:ARG:HB3	1.66	0.76
1:B:330:LYS:HZ3	1:B:572:ARG:HB3	1.50	0.76
1:A:118:GLN:HG2	1:A:130:LEU:HB2	1.66	0.75
1:A:483:HIS:H	1:B:448:GLN:NE2	1.85	0.74
1:B:539:LYS:HA	1:B:542:GLN:HE21	1.51	0.73
1:B:136:GLN:HA	1:B:139:GLN:HE21	1.53	0.73
1:A:136:GLN:HA	1:A:139:GLN:HE21	1.54	0.72
1:B:445:SER:HB3	1:B:449:MET:CE	2.20	0.71
1:A:448:GLN:NE2	1:B:483:HIS:H	1.89	0.71
1:B:141:LEU:HG	1:B:145:MET:HE1	1.73	0.70
1:A:141:LEU:HG	1:A:145:MET:HE2	1.74	0.69
1:A:445:SER:HB3	1:A:449:MET:CE	2.21	0.69
1:A:225:LYS:HG3	1:A:325:GLU:HG2	1.73	0.69
1:A:123:GLU:OE2	1:B:116:HIS:CD2	2.44	0.69
1:B:382:THR:HG23	2:B:611:TDP:O21	1.94	0.68
1:B:330:LYS:NZ	1:B:572:ARG:HD3	2.09	0.67
1:A:443:PHE:HE2	1:A:491:ALA:HA	1.62	0.65
1:B:237:LYS:HE3	1:B:388:ALA:HA	1.78	0.65
1:A:237:LYS:HE3	1:A:388:ALA:HA	1.79	0.64
1:A:133:SER:O	1:A:136:GLN:HG2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:VAL:HG11	1:A:307:ALA:HB2	1.79	0.64
1:A:371:ALA:N	1:A:426:GLN:HE22	1.95	0.64
1:B:443:PHE:HE2	1:B:491:ALA:HA	1.62	0.63
1:A:184:VAL:HG11	1:A:293:ILE:HD11	1.80	0.63
1:B:52:VAL:HG13	1:B:413:PRO:HB3	1.80	0.62
1:B:330:LYS:NZ	1:B:572:ARG:HB3	2.15	0.62
1:B:133:SER:O	1:B:136:GLN:HG2	1.99	0.61
1:B:184:VAL:HG11	1:B:293:ILE:HD11	1.81	0.61
2:A:611:TDP:H4A3	1:B:25:THR:O	2.00	0.61
1:A:137:ILE:HG22	1:A:138:PRO:HD3	1.82	0.61
1:A:52:VAL:HG13	1:A:413:PRO:HB3	1.83	0.61
1:A:312:ILE:HG12	3:A:612:FAD:C2A	2.31	0.60
1:A:383:PRO:HD3	2:A:611:TDP:O22	2.01	0.60
1:A:116:HIS:CD2	1:B:123:GLU:OE2	2.51	0.60
1:A:352:PRO:HD3	1:A:361:TYR:CE1	2.37	0.59
1:B:63:LEU:HD12	1:B:418:ALA:HA	1.84	0.59
1:B:304:VAL:HG11	1:B:307:ALA:HB2	1.84	0.59
1:A:447:VAL:HG23	1:A:496:ILE:HD11	1.85	0.58
1:B:547:SER:N	1:B:550:MET:HE3	2.15	0.58
1:B:371:ALA:N	1:B:426:GLN:HE22	2.00	0.58
1:B:290:GLN:HE22	1:B:299:GLY:H	1.52	0.58
1:B:3:GLN:HE22	1:B:8:TYR:HA	1.68	0.58
1:B:137:ILE:HG22	1:B:138:PRO:HD3	1.84	0.58
1:B:352:PRO:HD3	1:B:361:TYR:CE1	2.39	0.58
1:A:469:GLU:HG3	1:A:539:LYS:CE	2.34	0.58
1:B:467:ALA:C	1:B:469:GLU:H	2.08	0.57
1:B:447:VAL:HG23	1:B:496:ILE:HD11	1.85	0.57
1:A:137:ILE:CG2	1:A:138:PRO:HD3	2.34	0.57
1:B:367:SER:HB2	1:B:391:LEU:HD23	1.86	0.57
1:A:367:SER:HB2	1:A:391:LEU:HD23	1.87	0.57
1:B:118:GLN:CG	1:B:130:LEU:HB2	2.34	0.56
1:A:483:HIS:H	1:B:448:GLN:HE21	1.52	0.56
1:B:534:ILE:O	1:B:539:LYS:HE3	2.05	0.56
1:B:137:ILE:CG2	1:B:138:PRO:HD3	2.35	0.56
1:B:469:GLU:HG3	1:B:539:LYS:CE	2.35	0.56
1:B:136:GLN:HA	1:B:139:GLN:NE2	2.21	0.55
1:A:290:GLN:HE22	1:A:299:GLY:H	1.54	0.55
1:A:63:LEU:HD12	1:A:418:ALA:HA	1.89	0.55
1:B:248:VAL:O	1:B:248:VAL:HG12	2.06	0.55
1:B:251:THR:HG21	1:B:260:PHE:HB2	1.89	0.55
1:A:279:ARG:NH1	1:B:106:GLU:OE2	2.38	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:THR:O	2:B:611:TDP:H4A3	2.07	0.54
1:A:409:ALA:HB2	1:A:438:MET:HB3	1.90	0.54
1:A:136:GLN:HA	1:A:139:GLN:NE2	2.21	0.54
1:A:251:THR:HG21	1:A:260:PHE:HB2	1.89	0.53
1:A:468:MET:O	1:A:534:ILE:HD12	2.08	0.53
1:B:496:ILE:HG23	1:B:520:PRO:HB2	1.90	0.53
1:A:441:GLY:HA3	1:B:437:SER:O	2.09	0.53
1:A:443:PHE:CE2	1:A:491:ALA:HA	2.44	0.53
1:A:345:GLY:HA2	1:A:348:ASP:OD1	2.09	0.53
1:A:545:GLY:HA2	1:B:165:LEU:HD12	1.90	0.53
1:A:308:LEU:N	1:A:308:LEU:HD12	2.25	0.52
1:A:229:PRO:HB2	1:A:248:VAL:HG21	1.91	0.52
1:A:106:GLU:OE2	1:B:279:ARG:NH1	2.42	0.52
1:B:382:THR:N	1:B:383:PRO:CD	2.72	0.52
1:A:534:ILE:O	1:A:539:LYS:HE3	2.09	0.52
1:B:312:ILE:HG12	3:B:612:FAD:C2A	2.40	0.52
1:A:382:THR:N	1:A:383:PRO:CD	2.73	0.51
1:A:437:SER:O	1:B:441:GLY:HA3	2.10	0.51
1:B:443:PHE:CE2	1:B:491:ALA:HA	2.45	0.51
1:A:496:ILE:HG23	1:A:520:PRO:HB2	1.92	0.51
1:B:229:PRO:HB2	1:B:248:VAL:HG21	1.93	0.51
1:B:383:PRO:HD3	2:B:611:TDP:O22	2.11	0.51
1:A:446:VAL:HG22	1:A:496:ILE:HD11	1.92	0.50
1:A:248:VAL:O	1:A:248:VAL:HG12	2.10	0.50
1:B:409:ALA:HB2	1:B:438:MET:HB3	1.92	0.50
1:A:547:SER:H	1:A:550:MET:CE	2.25	0.50
1:A:118:GLN:CG	1:A:130:LEU:HB2	2.38	0.49
1:A:371:ALA:HB3	1:A:374:ALA:HB2	1.94	0.49
2:A:611:TDP:H352	4:B:621:SO4:O1	2.12	0.49
1:B:469:GLU:HG3	1:B:539:LYS:NZ	2.26	0.49
1:B:345:GLY:HA2	1:B:348:ASP:OD1	2.12	0.49
1:A:186:PRO:HG2	1:A:191:LEU:HD21	1.94	0.49
1:B:412:MET:HB3	1:B:413:PRO:HD3	1.95	0.49
1:B:446:VAL:HG22	1:B:496:ILE:HD11	1.94	0.49
1:A:469:GLU:HG3	1:A:539:LYS:NZ	2.27	0.49
1:B:241:GLU:HG3	1:B:241:GLU:O	2.12	0.48
1:A:460:ASN:CG	1:A:462:VAL:HG12	2.34	0.48
1:A:448:GLN:HE21	1:B:483:HIS:H	1.60	0.48
1:B:307:ALA:C	1:B:308:LEU:HD12	2.34	0.48
1:A:379:ASP:N	1:A:411:ALA:HB2	2.29	0.48
1:A:412:MET:HB3	1:A:413:PRO:HD3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:ARG:HD3	1:B:106:GLU:OE1	2.14	0.48
1:A:227:LYS:HD3	1:A:326:LYS:O	2.13	0.48
1:B:460:ASN:CG	1:B:462:VAL:HG12	2.34	0.48
1:B:197:LEU:HD22	1:B:289:ILE:HD11	1.96	0.48
3:A:612:FAD:H9	3:A:612:FAD:H1'1	1.59	0.47
1:A:293:ILE:O	1:A:293:ILE:HD13	2.15	0.47
1:B:522:LEU:HD23	1:B:522:LEU:C	2.35	0.47
1:B:227:LYS:N	1:B:227:LYS:HD2	2.30	0.47
1:A:467:ALA:HB2	1:A:471:LYS:HA	1.97	0.46
1:B:227:LYS:HD3	1:B:326:LYS:O	2.16	0.46
1:A:197:LEU:HD22	1:A:289:ILE:HD11	1.97	0.46
1:B:8:TYR:CE1	1:B:172:ALA:HB1	2.51	0.46
1:A:467:ALA:HB1	1:A:470:MET:C	2.36	0.46
1:B:467:ALA:HB1	1:B:470:MET:C	2.36	0.46
1:A:522:LEU:C	1:A:522:LEU:HD23	2.35	0.46
1:B:50:GLU:HB2	1:B:80:HIS:HB3	1.98	0.46
1:B:468:MET:O	1:B:534:ILE:HD12	2.16	0.45
1:B:437:SER:HA	1:B:440:MET:HB2	1.97	0.45
1:A:241:GLU:HG3	1:A:241:GLU:O	2.15	0.45
1:A:165:LEU:HD12	1:B:545:GLY:HA2	1.99	0.45
1:A:34:ASP:O	1:A:38:ARG:HG3	2.17	0.45
1:A:50:GLU:HB2	1:A:80:HIS:HB3	1.99	0.45
1:B:74:CYS:HA	1:B:100:ALA:O	2.17	0.45
1:B:464:GLY:HA2	2:B:611:TDP:O21	2.17	0.45
1:B:186:PRO:HG2	1:B:191:LEU:HD21	1.98	0.45
2:A:611:TDP:H4'2	2:A:611:TDP:H2	1.81	0.45
1:B:293:ILE:HD13	1:B:293:ILE:O	2.17	0.45
1:A:367:SER:HB2	1:A:391:LEU:HA	1.98	0.45
1:B:371:ALA:HB3	1:B:374:ALA:HB2	1.99	0.44
1:A:227:LYS:N	1:A:227:LYS:HD2	2.32	0.44
1:A:442:ASP:O	1:A:445:SER:HB2	2.18	0.44
1:A:361:TYR:O	1:A:365:GLN:HG2	2.18	0.44
1:B:379:ASP:N	1:B:411:ALA:HB2	2.32	0.44
1:A:75:GLY:HA3	1:A:76:PRO:HD3	1.83	0.44
1:B:320:LEU:HB2	1:B:321:PRO:HD3	1.99	0.44
1:A:359:PRO:HG2	1:A:530:GLU:HG3	2.00	0.44
1:B:308:LEU:HD12	1:B:308:LEU:N	2.32	0.44
1:A:547:SER:H	1:A:550:MET:HE2	1.83	0.43
1:B:467:ALA:HB2	1:B:471:LYS:HA	1.99	0.43
1:A:117:PRO:HB2	1:A:158:VAL:CG2	2.49	0.43
1:B:367:SER:HB2	1:B:391:LEU:HA	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:CYS:HA	1:A:100:ALA:O	2.18	0.43
4:A:621:SO4:O2	2:B:611:TDP:H352	2.18	0.43
1:A:320:LEU:HB2	1:A:321:PRO:HD3	2.01	0.43
1:A:339:TYR:O	1:A:343:ARG:HG3	2.19	0.43
1:A:348:ASP:O	1:A:351:LYS:HE3	2.18	0.42
1:A:68:ALA:O	1:A:95:VAL:HA	2.19	0.42
1:A:437:SER:HA	1:A:440:MET:HB2	2.00	0.42
1:A:464:GLY:HA2	2:A:611:TDP:O21	2.20	0.42
1:B:179:ALA:HA	1:B:180:PRO:HD2	1.94	0.42
1:B:258:SER:HB3	1:B:339:TYR:HA	2.01	0.42
1:B:359:PRO:HG2	1:B:530:GLU:HG3	2.02	0.42
2:B:611:TDP:C6'	2:B:611:TDP:H4A2	2.50	0.42
1:A:513:ARG:HD2	1:A:517:ILE:HD11	2.01	0.42
1:B:253:LEU:HD22	1:B:548:LEU:HB2	2.01	0.42
3:B:612:FAD:H9	3:B:612:FAD:H1'1	1.70	0.42
1:B:242:TYR:HB2	1:B:394:ASN:HA	2.01	0.42
1:A:469:GLU:OE1	1:A:539:LYS:HD3	2.20	0.41
1:B:290:GLN:NE2	1:B:299:GLY:H	2.16	0.41
1:A:513:ARG:O	1:A:517:ILE:HG13	2.20	0.41
1:A:126:HIS:CE1	1:A:153:GLY:HA3	2.55	0.41
1:A:253:LEU:HD22	1:A:548:LEU:HB2	2.01	0.41
1:B:137:ILE:N	1:B:138:PRO:CD	2.83	0.41
1:B:352:PRO:HD3	1:B:361:TYR:CZ	2.56	0.41
1:A:150:LEU:HD23	1:A:150:LEU:HA	1.83	0.41
1:A:251:THR:HG23	1:A:256:PHE:O	2.20	0.41
1:B:50:GLU:HB2	1:B:80:HIS:CB	2.51	0.41
1:B:540:LEU:HD23	1:B:540:LEU:HA	1.75	0.41
1:A:106:GLU:OE1	1:B:279:ARG:HD3	2.21	0.41
1:B:251:THR:HG23	1:B:256:PHE:O	2.19	0.41
1:B:126:HIS:CE1	1:B:153:GLY:HA3	2.56	0.41
1:B:112:PHE:O	1:B:113:GLN:HB2	2.21	0.41
1:B:467:ALA:C	1:B:469:GLU:N	2.73	0.41
1:A:137:ILE:N	1:A:138:PRO:CD	2.84	0.41
1:B:311:ASP:HA	3:B:612:FAD:N1A	2.36	0.41
1:B:469:GLU:OE1	1:B:539:LYS:HD3	2.21	0.40
1:B:117:PRO:HB2	1:B:158:VAL:HG21	2.03	0.40
1:B:86:PHE:O	1:B:90:ARG:HG2	2.20	0.40
1:B:442:ASP:O	1:B:445:SER:HB2	2.20	0.40
1:B:477:THR:C	1:B:479:GLY:N	2.75	0.40
1:A:347:ASP:O	1:A:348:ASP:C	2.59	0.40
1:B:276:PHE:HA	1:B:277:PRO:HD3	1.91	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:PRO:HD3	1:A:361:TYR:CZ	2.56	0.40
1:A:242:TYR:HB2	1:A:394:ASN:HA	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	569/572 (100%)	537 (94%)	30 (5%)	2 (0%)	39	74
1	B	569/572 (100%)	537 (94%)	30 (5%)	2 (0%)	39	74
All	All	1138/1144 (100%)	1074 (94%)	60 (5%)	4 (0%)	39	74

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	468	MET
1	A	469	GLU
1	B	469	GLU
1	A	467	ALA

### 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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	457/458 (100%)	437 (96%)	20 (4%)	35	70
1	B	457/458 (100%)	437 (96%)	20 (4%)	35	70
All	All	914/916 (100%)	874 (96%)	40 (4%)	35	70

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	THR
1	A	36	LEU
1	A	46	SER
1	A	66	GLU
1	A	118	GLN
1	A	139	GLN
1	A	146	ARG
1	A	173	THR
1	A	185	THR
1	A	187	GLU
1	A	210	SER
1	A	231	VAL
1	A	241	GLU
1	A	251	THR
1	A	284	THR
1	A	293	ILE
1	A	347	ASP
1	A	360	GLN
1	A	446	VAL
1	A	508	ASP
1	B	2	LYS
1	B	3	GLN
1	B	25	THR
1	B	36	LEU
1	B	46	SER
1	B	118	GLN
1	B	139	GLN
1	B	146	ARG
1	B	173	THR
1	B	185	THR
1	B	187	GLU
1	B	231	VAL

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Mol	Chain	Res	Type
1	B	241	GLU
1	B	251	THR
1	B	284	THR
1	B	293	ILE
1	B	347	ASP
1	B	360	GLN
1	B	446	VAL
1	B	508	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	92	HIS
1	A	116	HIS
1	A	126	HIS
1	A	139	GLN
1	A	151	ASN
1	A	181	GLN
1	A	290	GLN
1	A	426	GLN
1	A	448	GLN
1	A	542	GLN
1	B	3	GLN
1	B	92	HIS
1	B	116	HIS
1	B	126	HIS
1	B	139	GLN
1	B	151	ASN
1	B	181	GLN
1	B	290	GLN
1	B	426	GLN
1	B	448	GLN
1	B	542	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.



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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 2 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$
2	TDP	A	611	5	21,27,27	1.45	1 (4%)	31,40,40	1.88	9 (29%)
3	FAD	A	612	-	48,58,58	1.04	2 (4%)	54,89,89	2.06	8 (14%)
4	SO4	A	621	-	4,4,4	0.22	0	6,6,6	0.27	0
2	TDP	B	611	5	21,27,27	1.35	1 (4%)	31,40,40	2.03	9 (29%)
3	FAD	B	612	-	48,58,58	0.98	3 (6%)	54,89,89	2.08	8 (14%)
4	SO4	B	621	-	4,4,4	0.17	0	6,6,6	0.28	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
2	TDP	A	611	5	-	0/16/17/17	0/2/2/2
3	FAD	A	612	-	-	0/30/50/50	0/6/6/6
4	SO4	A	621	-	-	0/0/0/0	0/0/0/0
2	TDP	B	611	5	-	0/16/17/17	0/2/2/2
3	FAD	B	612	-	-	0/30/50/50	0/6/6/6
4	SO4	B	621	-	-	0/0/0/0	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	611	TDP	C4-N3	-5.91	1.34	1.39
2	B	611	TDP	C4-N3	-5.37	1.35	1.39
3	B	612	FAD	C10-N10	-3.72	1.34	1.39
3	A	612	FAD	C10-N10	-3.55	1.35	1.39
3	A	612	FAD	C9A-N10	-3.50	1.33	1.38
3	B	612	FAD	C9A-N10	-2.35	1.35	1.38
3	B	612	FAD	O4B-C1B	2.39	1.44	1.41

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	612	FAD	N3A-C2A-N1A	-10.29	121.01	128.89
3	A	612	FAD	N3A-C2A-N1A	-9.69	121.48	128.89
2	B	611	TDP	C5'-C35-N3	-4.29	106.15	113.33
2	B	611	TDP	C5A-C5-S1	-4.03	114.60	120.24
2	A	611	TDP	C5'-C35-N3	-3.91	106.80	113.33
3	A	612	FAD	C4X-C4-N3	-3.31	119.06	123.59
3	B	612	FAD	C4X-C4-N3	-3.24	119.16	123.59
2	A	611	TDP	C5A-C5-S1	-3.22	115.73	120.24
3	A	612	FAD	O3P-P-O5'	-3.14	94.60	102.94
3	B	612	FAD	C4A-C5A-N7A	-2.68	107.01	109.48
2	A	611	TDP	C5'-C6'-N1'	-2.64	119.28	123.86
2	B	611	TDP	C4A-C4-C5	-2.62	123.01	128.90
2	B	611	TDP	C5'-C6'-N1'	-2.48	119.56	123.86
3	B	612	FAD	P-O3P-PA	-2.46	125.83	132.73
2	B	611	TDP	N1'-C2'-N3'	-2.37	121.22	125.60
2	A	611	TDP	N1'-C2'-N3'	-2.31	121.33	125.60
3	A	612	FAD	C4A-C5A-N7A	-2.16	107.50	109.48
3	A	612	FAD	C1B-N9A-C4A	-2.13	123.72	126.94
2	A	611	TDP	C4A-C4-C5	-2.05	124.30	128.90
3	B	612	FAD	C9A-C5X-N5	-2.04	119.34	122.36
3	B	612	FAD	O4B-C1B-N9A	2.11	112.53	108.10
2	A	611	TDP	N4'-C4'-N3'	2.19	120.13	116.95
2	B	611	TDP	C6'-C5'-C4'	2.34	119.08	115.72
3	A	612	FAD	C4-C4X-N5	2.40	121.63	118.72
3	B	612	FAD	C5X-C9A-N10	2.57	119.57	117.62
3	A	612	FAD	C5X-C9A-N10	2.66	119.64	117.62
2	A	611	TDP	C6'-C5'-C4'	2.66	119.54	115.72
2	B	611	TDP	C6'-N1'-C2'	2.90	120.84	115.77
2	A	611	TDP	C6'-N1'-C2'	2.96	120.94	115.77
2	B	611	TDP	C4A-C4-N3	3.01	126.61	122.59
2	A	611	TDP	C5A-C5-C4	4.54	131.63	127.56
2	B	611	TDP	C5A-C5-C4	5.40	132.41	127.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	612	FAD	C4-N3-C2	6.90	121.21	115.25
3	A	612	FAD	C4-N3-C2	7.51	121.74	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	611	TDP	6	0
3	A	612	FAD	2	0
4	A	621	SO4	1	0
2	B	611	TDP	7	0
3	B	612	FAD	3	0
4	B	621	SO4	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	571/572 (99%)	-0.42	15 (2%) 59 54	50, 77, 113, 227	0
1	B	571/572 (99%)	-0.24	14 (2%) 61 55	49, 75, 114, 221	0
All	All	1142/1144 (99%)	-0.33	29 (2%) 61 55	49, 76, 114, 227	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	469	GLU	13.9
1	B	473	GLY	13.2
1	B	468	MET	11.1
1	B	474	GLY	9.9
1	B	475	TYR	9.8
1	A	469	GLU	9.5
1	A	474	GLY	8.5
1	A	468	MET	7.8
1	B	476	LEU	6.5
1	A	475	TYR	6.5
1	A	473	GLY	6.2
1	A	476	LEU	4.6
1	A	472	ALA	4.1
1	A	478	ASP	4.0
1	A	571	LEU	3.5
1	A	470	MET	3.0
1	B	470	MET	3.0
1	A	572	ARG	3.0
1	B	435	GLY	2.9
1	B	478	ASP	2.5
1	B	567	LYS	2.5
1	A	435	GLY	2.4
1	B	432	GLY	2.3
1	A	471	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	472	ALA	2.2
1	A	477	THR	2.2
1	B	379	ASP	2.1
1	A	355	LYS	2.1
1	B	355	LYS	2.1

## 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 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(Å <sup>2</sup> )	Q<0.9
4	SO4	A	621	5/5	0.97	0.15	1.17	74,82,95,108	0
5	MG	A	616	1/1	0.91	0.22	0.91	57,57,57,57	0
5	MG	B	617	1/1	0.92	0.29	0.75	59,59,59,59	0
4	SO4	B	621	5/5	0.99	0.16	0.48	68,71,82,87	0
2	TDP	A	611	26/26	0.98	0.17	-0.09	46,60,79,82	0
2	TDP	B	611	26/26	0.97	0.18	-0.18	34,64,74,122	0
3	FAD	A	612	53/53	0.99	0.12	-0.59	46,65,79,93	0
3	FAD	B	612	53/53	0.98	0.12	-0.92	49,66,84,91	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.