



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 01:53 am GMT

PDB ID : 3LQ4
Title : E. coli pyruvate dehydrogenase complex E1 E235A mutant with high TDP concentration
Authors : Furey, W.
Deposited on : 2010-02-08
Resolution : 1.98 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

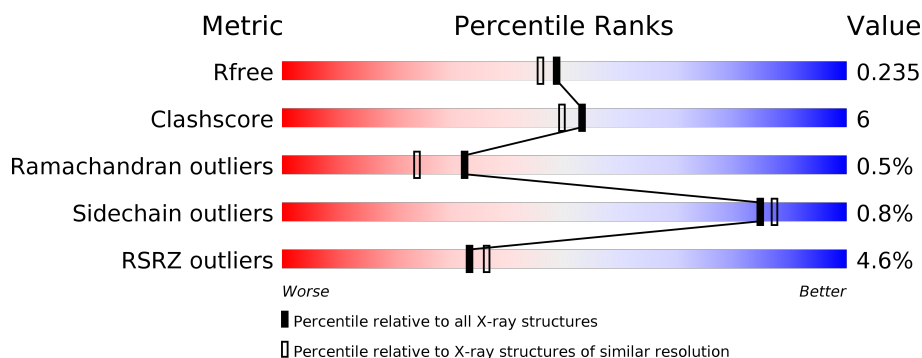
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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}	100719	9293 (2.00-1.96)
Clashscore	112137	10621 (2.00-1.96)
Ramachandran outliers	110173	10502 (2.00-1.96)
Sidechain outliers	110143	10501 (2.00-1.96)
RSRZ outliers	101464	9395 (2.00-1.96)

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 ≥ 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	886	<div> <div>5%</div> <div> <div></div> <div>77%</div> <div>13%</div> <div>10%</div> </div> </div>
1	B	886	<div> <div>3%</div> <div> <div></div> <div>81%</div> <div>9%</div> <div>10%</div> </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	MG	A	888	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13440 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.

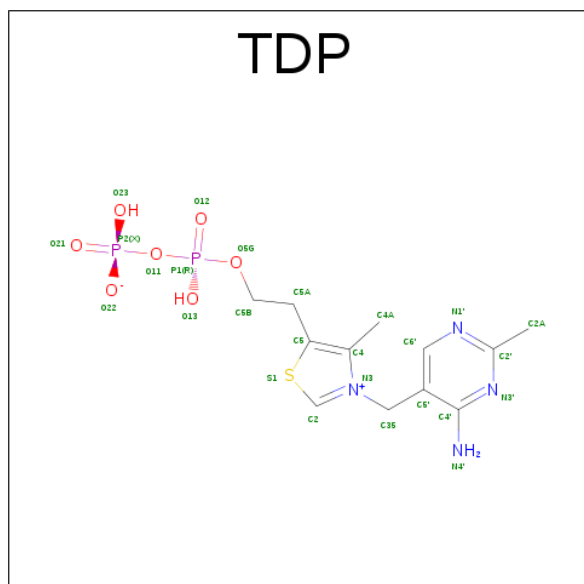
- Molecule 1 is a protein called Pyruvate dehydrogenase E1 component.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	801	Total	C	N	O	S	0	1	0
			6347	4022	1096	1203	26			
1	B	801	Total	C	N	O	S	0	0	0
			6337	4016	1093	1202	26			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	235	ALA	GLU	ENGINEERED	UNP P0AFG9
B	235	ALA	GLU	ENGINEERED	UNP P0AFG9

- 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		

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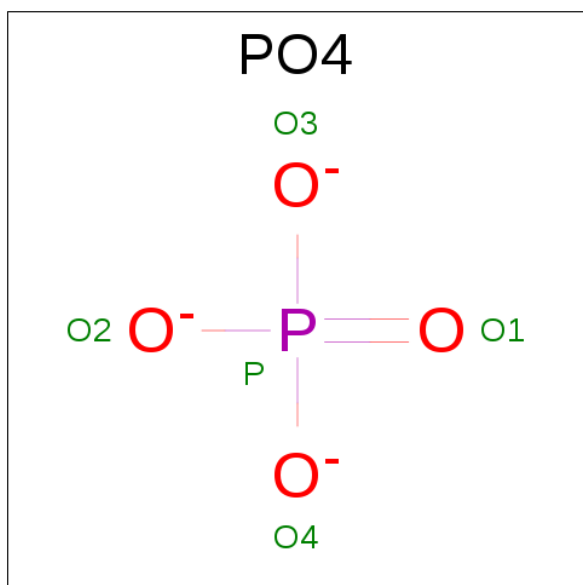
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	S	
			26	12	4	7	2	1	
								0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

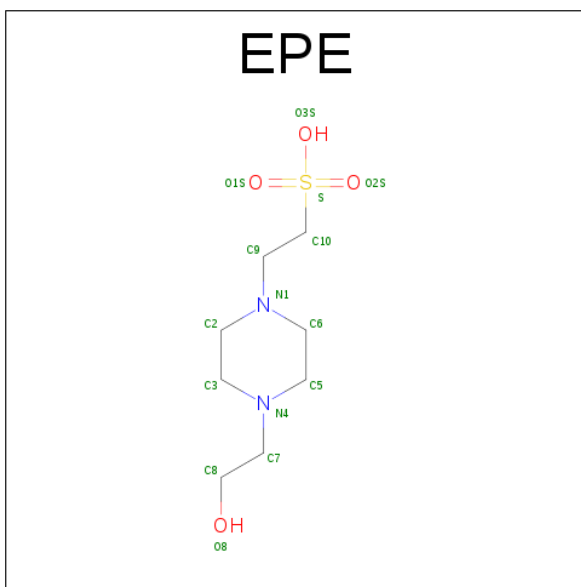
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg		
			1	1	0	0
3	A	1	Total	Mg		
			1	1	0	0

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

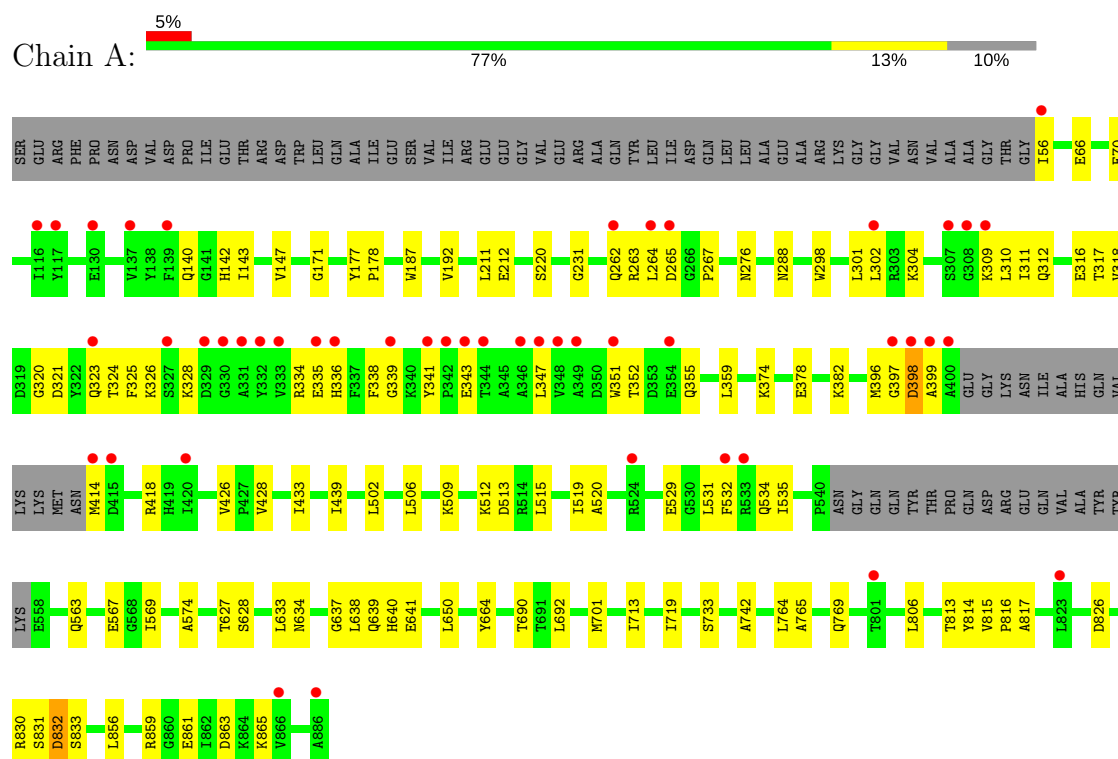
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	340	Total	O	0	0
			340	340		
6	B	337	Total	O	0	0
			337	337		

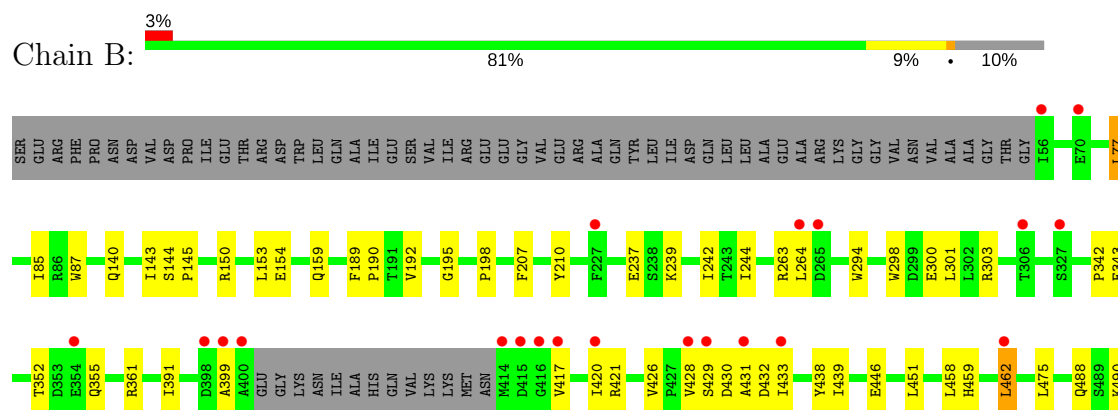
3 Residue-property plots

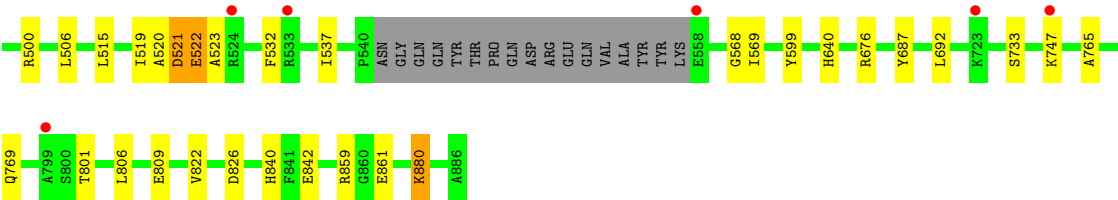
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Pyruvate dehydrogenase E1 component



• Molecule 1: Pyruvate dehydrogenase E1 component





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.21Å 142.22Å 82.85Å 90.00° 101.87° 90.00°	Depositor
Resolution (Å)	32.52 – 1.98 32.52 – 1.96	Depositor EDS
% Data completeness (in resolution range)	83.7 (32.52-1.98) 82.6 (32.52-1.96)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.06 (at 1.97Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.199 , 0.236 0.201 , 0.235	Depositor DCC
R_{free} test set	5434 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.366	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.022 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13440	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EPE, PO4, TDP, 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	0.32	0/6491	0.57	1/8776 (0.0%)
1	B	0.33	0/6480	0.58	0/8761
All	All	0.32	0/12971	0.57	1/17537 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	171	GLY	N-CA-C	5.22	126.15	113.10

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	6347	0	6184	92	0
1	B	6337	0	6178	72	0
2	A	26	0	16	7	0
2	B	26	0	16	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	15	0	18	3	0
6	A	340	0	0	2	0
6	B	337	0	0	4	0
All	All	13440	0	12412	155	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 6.

All (155) 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:887:TDP:C2	2:A:887:TDP:H2	0.97	1.49
2:B:887:TDP:H2	2:B:887:TDP:C2	0.97	1.48
1:A:302:LEU:HD11	1:A:311:ILE:HD11	1.51	0.93
1:B:421:ARG:HD3	1:B:433:ILE:HD11	1.53	0.90
1:A:192:VAL:HG21	1:B:640:HIS:HE1	1.35	0.89
1:B:421:ARG:CD	1:B:433:ILE:HD11	2.13	0.78
1:B:421:ARG:HD3	1:B:433:ILE:CD1	2.13	0.78
1:A:323:GLN:HA	1:A:326:LYS:HE3	1.65	0.76
1:B:264:LEU:HB2	2:B:887:TDP:H5A2	1.72	0.72
1:A:192:VAL:CG2	1:B:640:HIS:HE1	2.02	0.72
1:A:320:GLY:O	1:A:323:GLN:HG2	1.89	0.72
1:A:192:VAL:HG21	1:B:640:HIS:CE1	2.22	0.70
2:A:887:TDP:H5A1	1:B:569:ILE:HD11	1.76	0.67
1:A:264:LEU:HB2	2:A:887:TDP:H5A2	1.78	0.65
2:A:887:TDP:H4A1	1:B:569:ILE:HG12	1.80	0.63
1:B:154:GLU:HG2	1:B:451:LEU:HD21	1.81	0.63
1:A:640:HIS:HE1	1:B:192:VAL:HG21	1.63	0.63
1:A:426:VAL:CG1	1:A:439:ILE:HD11	2.29	0.62
1:A:832:ASP:OD2	1:A:833:SER:N	2.25	0.61
1:A:302:LEU:CD1	1:A:311:ILE:HD11	2.29	0.60
1:A:426:VAL:HG12	1:A:428:VAL:HG23	1.84	0.59
1:A:426:VAL:HG13	1:A:439:ILE:HD11	1.84	0.59
1:A:765:ALA:O	1:A:769:GLN:HG3	2.02	0.59
1:B:522:GLU:HG2	1:B:599:TYR:HE1	1.69	0.58
1:B:429:SER:HB3	1:B:432:ASP:OD2	2.04	0.58
1:A:640:HIS:HE1	1:B:192:VAL:CG2	2.16	0.57
1:A:192:VAL:CG2	1:B:640:HIS:CE1	2.86	0.57
1:B:809:GLU:HG2	1:B:822:VAL:HG21	1.86	0.57
1:A:863:ASP:OD2	1:A:865:LYS:HB2	2.04	0.57
1:A:352:THR:OG1	1:A:355:GLN:HG3	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:856:LEU:HD22	1:A:861:GLU:OE1	2.04	0.56
1:A:70:GLU:OE2	5:A:890:EPE:H21	2.05	0.56
1:B:522:GLU:HG2	1:B:599:TYR:CE1	2.41	0.56
1:A:140:GLN:O	1:A:143:ILE:HG13	2.06	0.56
1:B:417:VAL:O	1:B:420:ILE:HG12	2.06	0.55
1:A:304:LYS:HE3	1:A:347:LEU:HD23	1.88	0.55
1:B:801:THR:HB	6:B:1199:HOH:O	2.05	0.55
1:B:842:GLU:OE2	1:B:880:LYS:HE3	2.07	0.54
1:A:567:GLU:HG3	1:A:574:ALA:HA	1.89	0.54
1:A:813:THR:HG23	1:A:814:TYR:CE1	2.42	0.54
2:A:887:TDP:H2	6:B:1137:HOH:O	2.08	0.54
1:A:56:ILE:HD13	1:A:276:ASN:HB3	1.90	0.53
1:A:304:LYS:HE3	1:A:347:LEU:CD2	2.39	0.53
1:B:429:SER:C	1:B:431:ALA:H	2.11	0.53
1:B:77:LEU:HD11	1:B:446:GLU:HG2	1.89	0.53
1:A:813:THR:CG2	1:A:814:TYR:CZ	2.91	0.53
1:A:529:GLU:N	1:A:529:GLU:OE1	2.41	0.52
1:B:421:ARG:HD2	1:B:428:VAL:HG13	1.91	0.52
1:A:212:GLU:OE2	1:A:220:SER:HB3	2.08	0.52
1:A:263:ARG:HG3	1:A:264:LEU:H	1.75	0.52
1:A:529:GLU:HA	1:A:532:PHE:CD2	2.45	0.52
1:B:747:LYS:HG2	6:B:989:HOH:O	2.10	0.52
1:A:264:LEU:HD13	1:B:522:GLU:OE1	2.09	0.51
1:B:150:ARG:O	1:B:154:GLU:HG3	2.10	0.51
1:A:265:ASP:OD2	1:B:522:GLU:HA	2.10	0.51
1:B:140:GLN:O	1:B:143:ILE:HG13	2.11	0.51
1:A:414:MET:HE1	1:A:433:ILE:HG23	1.94	0.50
1:B:519:ILE:HD12	1:B:523:ALA:HB2	1.93	0.50
1:B:237:GLU:H	1:B:237:GLU:CD	2.15	0.50
1:B:506:LEU:HD23	1:B:515:LEU:HD12	1.93	0.50
1:A:831:SER:O	1:A:832:ASP:HB2	2.10	0.50
2:A:887:TDP:C4A	1:B:569:ILE:HG12	2.42	0.50
1:B:85:ILE:HG12	1:B:153:LEU:HD12	1.93	0.50
1:B:859:ARG:NH1	1:B:861:GLU:CD	2.65	0.49
1:B:488:GLN:OE1	1:B:500:ARG:NH1	2.45	0.49
1:A:531:LEU:HD23	1:A:534:GLN:HE21	1.77	0.49
1:A:719:ILE:HD12	1:A:742:ALA:HB1	1.93	0.49
1:A:813:THR:HG23	1:A:814:TYR:CD1	2.47	0.49
1:A:312:GLN:O	1:A:316:GLU:HG2	2.12	0.49
1:A:414:MET:O	1:A:418:ARG:HG3	2.13	0.49
1:B:244:ILE:HD12	1:B:244:ILE:N	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142[O]:HIS:N	1:A:142[O]:HIS:ND1	2.58	0.48
1:A:262:GLN:HA	1:A:267:PRO:HA	1.96	0.48
1:A:374:LYS:O	1:A:378:GLU:HG3	2.13	0.48
1:A:288:ASN:HB2	1:A:382:LYS:HE3	1.96	0.48
1:B:532:PHE:CE2	1:B:537:ILE:HD11	2.48	0.48
1:A:262:GLN:HB2	6:A:912:HOH:O	2.13	0.48
1:B:144:SER:OG	1:B:145:PRO:HD3	2.14	0.48
1:A:806:LEU:HD11	1:B:806:LEU:HD11	1.96	0.48
1:A:309:LYS:HE3	1:A:341:TYR:CD2	2.49	0.47
1:A:512:LYS:HG3	1:A:513:ASP:N	2.29	0.47
1:A:713:ILE:HB	1:A:764:LEU:HD11	1.96	0.47
1:B:765:ALA:O	1:B:769:GLN:HG3	2.14	0.47
1:A:318:VAL:HG22	1:A:321:ASP:OD2	2.14	0.47
1:B:361:ARG:HD2	1:B:391:ILE:HG13	1.97	0.46
1:B:429:SER:C	1:B:431:ALA:N	2.68	0.46
1:A:634:ASN:HB2	1:A:832:ASP:O	2.15	0.46
1:B:159:GLN:HG3	1:B:438:TYR:CD2	2.50	0.46
1:B:87:TRP:CD1	1:B:439:ILE:HG12	2.50	0.46
1:A:426:VAL:HG13	1:A:439:ILE:CD1	2.46	0.46
1:B:521:ASP:HB2	1:B:568:GLY:HA2	1.97	0.46
1:B:352:THR:OG1	1:B:355:GLN:HG3	2.16	0.46
1:A:325:PHE:HE1	1:A:336:HIS:HB2	1.81	0.46
1:B:207:PHE:O	1:B:210:TYR:HB3	2.16	0.46
1:B:426:VAL:HG12	1:B:428:VAL:HG12	1.98	0.46
1:A:692:LEU:HD13	1:A:733:SER:HB3	1.97	0.45
1:B:430:ASP:O	1:B:430:ASP:OD1	2.33	0.45
1:A:231:GLY:C	1:B:569:ILE:HD12	2.37	0.45
1:B:300:GLU:HG3	6:B:1140:HOH:O	2.17	0.45
1:A:650:LEU:HD12	1:A:650:LEU:C	2.37	0.45
1:B:859:ARG:HH12	1:B:861:GLU:CD	2.20	0.45
1:A:142[A]:HIS:NE2	2:A:887:TDP:S1	2.89	0.45
1:A:301:LEU:HD22	1:A:310:LEU:HD22	1.97	0.45
1:A:502:LEU:HD23	1:A:531:LEU:HD11	1.99	0.45
1:A:519:ILE:HG22	1:A:520:ALA:N	2.32	0.45
1:A:66:GLU:O	5:A:890:EPE:H81	2.17	0.45
1:B:189:PHE:HA	1:B:190:PRO:HD3	1.78	0.45
1:A:396:MET:O	1:A:398:ASP:N	2.50	0.44
1:B:264:LEU:O	1:B:264:LEU:HD23	2.17	0.44
1:A:192:VAL:HG22	1:A:192:VAL:O	2.17	0.44
1:A:506:LEU:HD23	1:A:515:LEU:HD12	2.00	0.44
1:A:531:LEU:HD23	1:A:534:GLN:NE2	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:639:GLN:OE1	1:B:192:VAL:HG12	2.18	0.43
1:B:303:ARG:HH11	1:B:303:ARG:HG3	1.83	0.43
1:B:676:ARG:HD3	1:B:687:TYR:OH	2.18	0.43
1:A:343:GLU:HA	6:A:956:HOH:O	2.17	0.43
1:A:664:TYR:CG	1:A:701:MET:HB2	2.53	0.43
1:A:815:VAL:HG12	1:A:817:ALA:H	1.83	0.43
1:A:637:GLY:O	1:A:641:GLU:HG3	2.18	0.43
1:A:351:TRP:HA	1:A:355:GLN:OE1	2.19	0.43
1:B:520:ALA:O	1:B:522:GLU:N	2.52	0.43
1:A:264:LEU:HB3	1:B:521:ASP:OD2	2.19	0.43
1:B:421:ARG:CD	1:B:428:VAL:HG13	2.49	0.43
1:B:342:PRO:HG2	1:B:343:GLU:OE1	2.19	0.43
1:A:628:SER:OG	1:A:690:THR:HB	2.19	0.42
1:B:840:HIS:O	1:B:880:LYS:NZ	2.48	0.42
1:A:320:GLY:O	1:A:323:GLN:CG	2.64	0.42
1:B:462:LEU:C	1:B:462:LEU:HD23	2.40	0.42
1:B:458:LEU:O	1:B:459:HIS:HB2	2.19	0.42
1:A:177:TYR:HB3	1:A:178:PRO:CD	2.50	0.42
1:A:569:ILE:HD11	2:B:887:TDP:H5A1	2.02	0.42
1:A:815:VAL:HA	1:A:816:PRO:HD3	1.86	0.42
1:A:535:ILE:HB	1:A:563:GLN:HB3	2.02	0.41
1:B:692:LEU:HD13	1:B:733:SER:HB3	2.01	0.41
1:A:178:PRO:HA	1:A:187:TRP:CG	2.55	0.41
1:A:317:THR:CG2	1:A:321:ASP:HB2	2.51	0.41
1:A:321:ASP:O	1:A:324:THR:HB	2.20	0.41
1:A:335:GLU:O	1:A:339:GLY:HA3	2.20	0.41
1:B:87:TRP:CD2	1:B:426:VAL:HG11	2.56	0.41
1:B:490:LYS:HE3	1:B:490:LYS:HB2	1.93	0.41
1:A:638:LEU:C	1:A:638:LEU:HD23	2.40	0.41
1:A:509:LYS:HA	1:A:512:LYS:HE3	2.03	0.41
1:B:195:GLY:O	1:B:198:PRO:HG2	2.21	0.41
1:A:326:LYS:HG2	1:A:326:LYS:O	2.20	0.41
1:A:298:TRP:CE2	1:A:359:LEU:HB3	2.56	0.41
1:A:813:THR:CG2	1:A:814:TYR:CE2	3.04	0.41
1:A:70:GLU:HG2	5:A:890:EPE:H21	2.02	0.41
1:B:264:LEU:HD23	1:B:264:LEU:C	2.40	0.41
1:A:627:THR:HB	1:A:633:LEU:HD13	2.03	0.40
1:A:859:ARG:CB	1:A:859:ARG:HH11	2.34	0.40
1:B:294:TRP:HB3	1:B:298:TRP:CD1	2.57	0.40
1:B:421:ARG:HD3	1:B:433:ILE:HD12	2.00	0.40
1:A:143:ILE:HD12	1:A:147:VAL:HG23	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:ARG:HA	1:A:338:PHE:HB2	2.02	0.40
1:B:239:LYS:O	1:B:242:ILE:HG12	2.21	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	796/886 (90%)	756 (95%)	35 (4%)	5 (1%)	28	20
1	B	795/886 (90%)	759 (96%)	33 (4%)	3 (0%)	38	30
All	All	1591/1772 (90%)	1515 (95%)	68 (4%)	8 (0%)	32	24

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	397	GLY
1	A	328	LYS
1	A	399	ALA
1	B	399	ALA
1	A	398	ASP
1	B	263	ARG
1	B	521	ASP
1	A	832	ASP

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	665/734 (91%)	662 (100%)	3 (0%)	91	93
1	B	664/734 (90%)	657 (99%)	7 (1%)	78	80
All	All	1329/1468 (90%)	1319 (99%)	10 (1%)	85	87

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	211	LEU
1	A	826	ASP
1	A	830	ARG
1	B	77	LEU
1	B	301	LEU
1	B	462	LEU
1	B	475	LEU
1	B	522	GLU
1	B	826	ASP
1	B	880	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	640	HIS

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 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	887	3	21,27,27	1.56	5 (23%)	25,40,40	1.14	2 (8%)
4	PO4	A	889	-	4,4,4	0.77	0	6,6,6	0.70	0
5	EPE	A	890	-	15,15,15	1.33	1 (6%)	18,20,20	2.42	5 (27%)
2	TDP	B	887	3	21,27,27	1.54	4 (19%)	25,40,40	1.12	2 (8%)
4	PO4	B	889	-	4,4,4	0.84	0	6,6,6	0.70	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	887	3	-	0/16/17/17	0/2/2/2
4	PO4	A	889	-	-	0/0/0/0	0/0/0/0
5	EPE	A	890	-	-	0/9/19/19	0/1/1/1
2	TDP	B	887	3	-	0/16/17/17	0/2/2/2
4	PO4	B	889	-	-	0/0/0/0	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	887	TDP	C6'-N1'	2.03	1.38	1.34
2	B	887	TDP	C2'-N1'	2.43	1.38	1.34
2	A	887	TDP	C5'-C4'	2.46	1.48	1.42
2	A	887	TDP	C2'-N1'	2.52	1.38	1.34
2	A	887	TDP	C4-N3	2.55	1.41	1.39
2	B	887	TDP	C4-N3	2.59	1.42	1.39
2	B	887	TDP	C5'-C4'	2.61	1.49	1.42
2	B	887	TDP	C4'-N3'	3.22	1.39	1.35
2	A	887	TDP	C4'-N3'	3.60	1.40	1.35
5	A	890	EPE	C10-S	3.97	1.83	1.77

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	890	EPE	O3S-S-O1S	-3.89	102.46	111.37
5	A	890	EPE	O3S-S-O2S	-3.78	102.70	111.37
2	B	887	TDP	O23-P2-O22	2.26	116.73	107.61
5	A	890	EPE	O2S-S-O1S	2.32	121.89	113.86
2	A	887	TDP	O23-P2-O22	2.40	117.29	107.61
2	B	887	TDP	C6'-N1'-C2'	2.45	120.12	115.88
2	A	887	TDP	C6'-N1'-C2'	2.53	120.25	115.88
5	A	890	EPE	O2S-S-C10	5.33	111.37	106.79
5	A	890	EPE	O1S-S-C10	5.47	111.49	106.79

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	887	TDP	7	0
5	A	890	EPE	3	0
2	B	887	TDP	3	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	801/886 (90%)	0.29	47 (5%)	23 26	22, 34, 72, 87	0
1	B	801/886 (90%)	0.11	27 (3%)	46 49	19, 33, 57, 87	0
All	All	1602/1772 (90%)	0.20	74 (4%)	33 36	19, 34, 64, 87	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	400	ALA	12.5
1	B	399	ALA	10.8
1	B	400	ALA	9.5
1	A	398	ASP	7.4
1	B	414	MET	6.7
1	A	399	ALA	6.6
1	A	397	GLY	6.4
1	A	331	ALA	5.2
1	A	335	GLU	5.1
1	A	342	PRO	4.9
1	A	302	LEU	4.6
1	A	348	VAL	4.5
1	A	329	ASP	4.3
1	B	306	THR	4.2
1	B	415	ASP	4.1
1	A	344	THR	4.0
1	B	416	GLY	3.9
1	A	349	ALA	3.9
1	A	323	GLN	3.7
1	A	264	LEU	3.7
1	A	341	TYR	3.4
1	B	398	ASP	3.4
1	B	420	ILE	3.4
1	A	330	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	354	GLU	3.4
1	A	332	TYR	3.4
1	B	428	VAL	3.3
1	A	117	TYR	3.3
1	A	265	ASP	3.1
1	A	347	LEU	3.1
1	A	354	GLU	3.0
1	A	351	TRP	3.0
1	B	417	VAL	3.0
1	A	533	ARG	2.9
1	A	886	ALA	2.9
1	B	524	ARG	2.9
1	B	56	ILE	2.9
1	B	265	ASP	2.8
1	A	339	GLY	2.8
1	A	420	ILE	2.8
1	B	264	LEU	2.7
1	A	415	ASP	2.7
1	B	533	ARG	2.7
1	B	462	LEU	2.6
1	A	309	LYS	2.6
1	A	56	ILE	2.5
1	B	431	ALA	2.5
1	A	137	VAL	2.5
1	A	336	HIS	2.5
1	A	866	VAL	2.5
1	A	343	GLU	2.4
1	B	327	SER	2.4
1	A	262	GLN	2.4
1	A	307	SER	2.3
1	A	532	PHE	2.3
1	A	327	SER	2.3
1	A	346	ALA	2.3
1	A	524	ARG	2.3
1	A	308	GLY	2.3
1	B	558	GLU	2.3
1	A	333	VAL	2.2
1	B	433	ILE	2.2
1	A	414	MET	2.2
1	A	801	THR	2.2
1	A	130	GLU	2.2
1	B	227	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	70	GLU	2.2
1	B	747	LYS	2.2
1	A	116	ILE	2.1
1	B	429	SER	2.1
1	B	799	ALA	2.1
1	B	723	LYS	2.1
1	A	823	LEU	2.1
1	A	139	PHE	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 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	MG	A	888	1/1	0.80	0.26	2.06	44,44,44,44	1
3	MG	B	888	1/1	0.89	0.17	0.82	36,36,36,36	0
5	EPE	A	890	15/15	0.88	0.14	0.60	43,46,62,63	0
2	TDP	A	887	26/26	0.94	0.14	-0.22	37,42,44,45	17
2	TDP	B	887	26/26	0.97	0.11	-0.31	34,40,41,43	7
4	PO4	A	889	5/5	0.98	0.07	-2.23	46,48,49,49	0
4	PO4	B	889	5/5	0.98	0.07	-2.87	57,58,58,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.