



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:42 PM UTC

PDB ID : 2G25 / pdb_00002g25
Title : E. Coli Pyruvate Dehydrogenase Phosphonolactylthiamin Diphosphate Complex
Authors : Furey, W.; Arjunan, P.; Chandrasekhar, K.
Deposited on : 2006-02-15
Resolution : 2.10 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

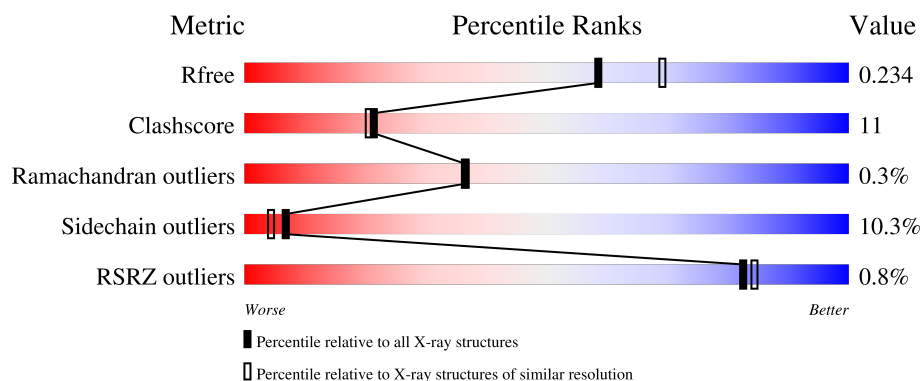
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	886	68% 22% • 6%
1	B	886	69% 21% • 6%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 13841 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	831	Total	C	N	O	S	0	0	0
			6591	4172	1140	1252	27			
1	B	831	Total	C	N	O	S	0	0	0
			6591	4172	1140	1252	27			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

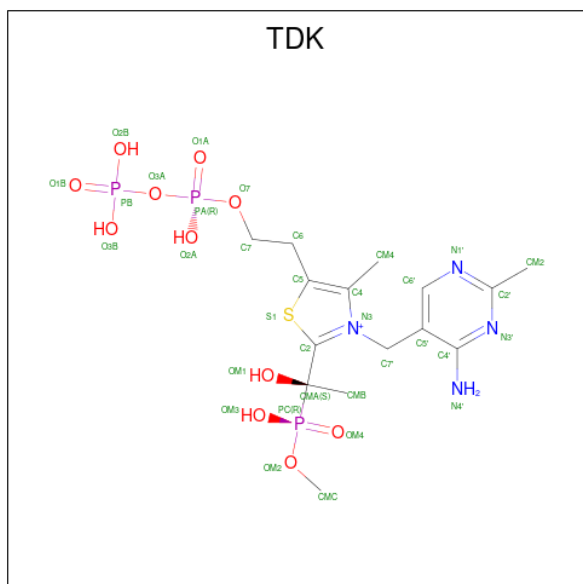
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 4 is 3-[(4-AMINO-2-METHYLPYRIMIDIN-5-YL)METHYL]-2-[(1S)-1-HYDROXY-1-[(R)-HYDROXY(METHOXY)PHOSPHORYL]ETHYL]-5-(2-[(S)-HYDROXY(PHOSPHONOOXY)PHOSPHORYL]OXY)ETHYL)-4-METHYL-1,3-THIAZOL-3-IUM (CCD ID: TDK) (formula: C₁₅H₂₆N₄O₁₁P₃S).



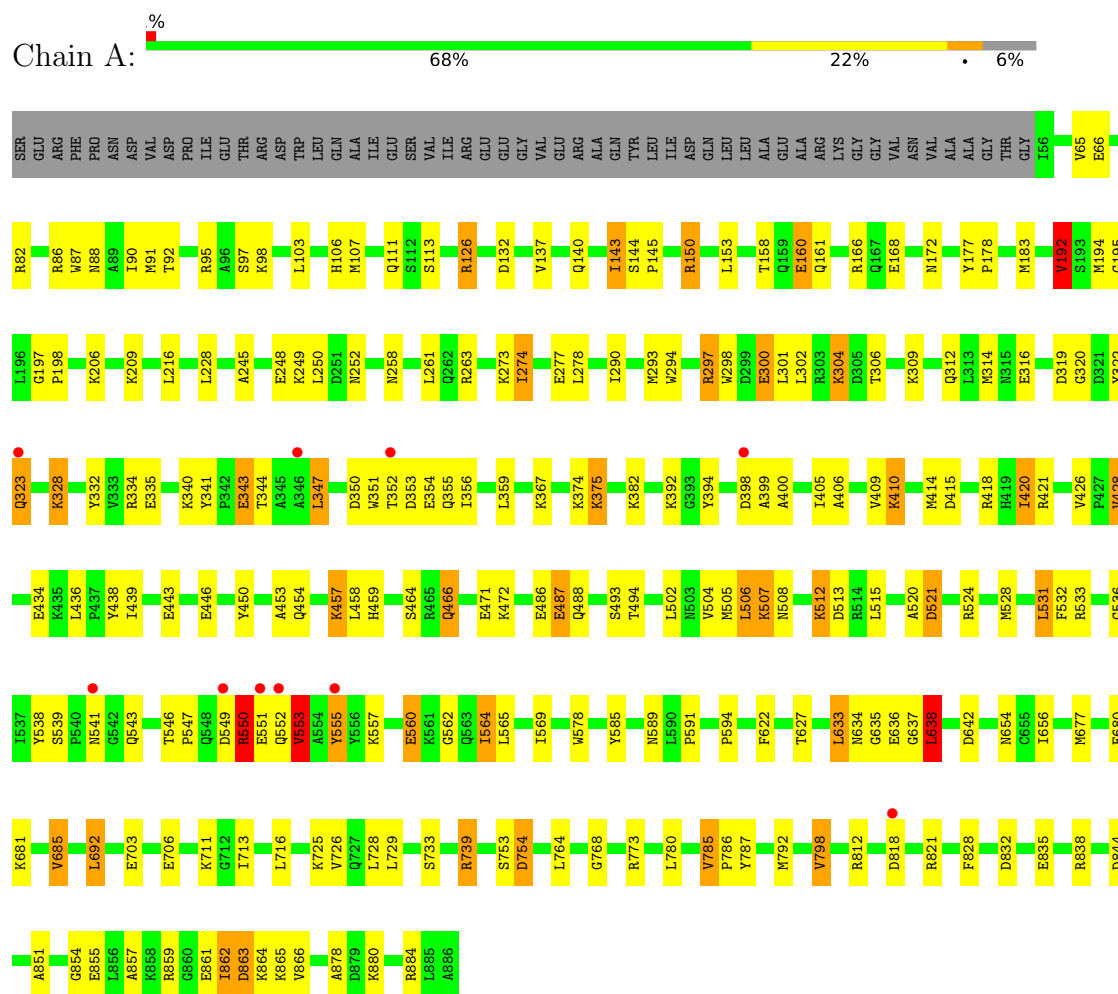
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	291	Total 291	O 291	0	0

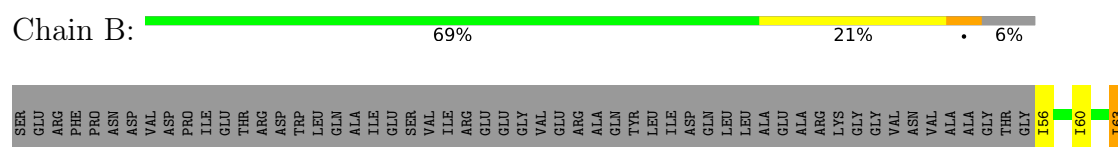
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: 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	81.57Å 142.51Å 82.09Å 90.00° 102.37° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10 8.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	94.0 (8.00-2.10) 91.9 (8.00-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.85 (at 2.10Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.185 , 0.244 0.180 , 0.234	Depositor DCC
R_{free} test set	5035 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	10.8	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 61.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.028 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13841	wwPDB-VP
Average B, all atoms (Å ²)	10.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, PO4, TDK

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	0.44	0/6741	0.96	25/9114 (0.3%)
1	B	0.44	0/6741	0.94	21/9114 (0.2%)
All	All	0.44	0/13482	0.95	46/18228 (0.3%)

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	B	0	2

There are no bond length outliers.

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	634	ASN	N-CA-C	10.07	123.51	111.33
1	B	634	ASN	N-CA-C	9.18	122.13	111.11
1	B	171	GLY	N-CA-C	8.49	125.07	114.37
1	B	633	LEU	N-CA-C	-8.04	100.52	110.65
1	B	636	GLU	N-CA-C	7.29	120.28	111.82
1	A	183	MET	CA-C-N	7.29	126.96	119.82
1	A	183	MET	C-N-CA	7.29	126.96	119.82
1	A	633	LEU	N-CA-C	-7.14	100.75	110.68
1	A	553	VAL	N-CA-C	6.80	123.48	109.34
1	A	552	GLN	N-CA-C	6.57	120.74	110.17
1	A	399	ALA	N-CA-C	6.55	118.97	111.11
1	A	785	VAL	N-CA-C	6.47	114.30	107.76
1	B	143	ILE	N-CA-C	-6.38	105.13	111.77
1	A	143	ILE	N-CA-C	-6.21	106.38	111.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	541	ASN	N-CA-C	-6.09	103.51	111.74
1	B	229	GLY	N-CA-C	-5.93	103.59	113.02
1	A	638	LEU	N-CA-C	5.82	120.91	111.37
1	A	637	GLY	N-CA-C	5.80	122.74	112.53
1	A	293	MET	N-CA-C	5.77	118.69	111.24
1	B	794	ASP	N-CA-C	5.71	119.94	112.92
1	A	636	GLU	N-CA-C	5.65	118.37	111.82
1	B	293	MET	N-CA-C	5.63	120.18	111.56
1	B	638	LEU	N-CA-C	5.48	120.42	111.37
1	A	550	ARG	N-CA-C	5.42	122.35	110.80
1	B	63	ILE	N-CA-C	5.34	112.84	107.55
1	A	521	ASP	N-CA-C	-5.29	99.52	110.80
1	B	577	SER	N-CA-C	-5.29	105.59	111.36
1	A	642	ASP	N-CA-C	5.27	116.89	108.41
1	A	504	VAL	N-CA-C	-5.26	105.37	110.42
1	A	137	VAL	N-CA-C	5.25	115.33	107.51
1	B	183	MET	CA-C-N	5.22	124.89	119.56
1	B	183	MET	C-N-CA	5.22	124.89	119.56
1	B	534	GLN	N-CA-C	5.19	116.62	111.07
1	A	192	VAL	N-CA-C	5.18	117.52	111.05
1	A	706	GLU	N-CA-C	5.18	117.32	111.11
1	A	415	ASP	N-CA-C	-5.17	105.72	111.36
1	B	679	GLY	N-CA-C	-5.17	105.67	112.81
1	B	199	ILE	CB-CA-C	-5.16	105.10	112.22
1	B	362	GLY	N-CA-C	5.15	118.87	112.64
1	A	754	ASP	N-CA-C	-5.11	100.19	108.52
1	A	466	GLN	CA-C-N	5.08	124.87	119.28
1	A	466	GLN	C-N-CA	5.08	124.87	119.28
1	B	652	ILE	CA-C-N	5.07	126.17	119.84
1	B	652	ILE	C-N-CA	5.07	126.17	119.84
1	B	460	GLY	N-CA-C	5.06	116.67	111.56
1	B	242	ILE	N-CA-C	5.04	117.35	111.05

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	598	TYR	Sidechain
1	B	686	TYR	Sidechain

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	6591	0	6422	149	0
1	B	6591	0	6422	143	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	10	0	0	0	0
3	B	15	0	0	2	0
4	A	34	0	22	2	0
4	B	34	0	22	4	0
5	A	273	0	0	7	0
5	B	291	0	0	4	0
All	All	13841	0	12888	284	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 11.

All (284) 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:550:ARG:HA	1:A:550:ARG:HH11	1.18	1.06
1:A:309:LYS:HG3	1:A:344:THR:HG23	1.52	0.91
1:A:549:ASP:HB3	1:A:555:TYR:HA	1.53	0.91
1:A:550:ARG:HH11	1:A:550:ARG:CA	1.87	0.88
1:A:550:ARG:HA	1:A:550:ARG:NH1	1.89	0.86
1:B:311:ILE:HA	1:B:314:MET:HE2	1.64	0.80
1:A:91:MET:O	1:A:95:ARG:HG3	1.84	0.77
1:B:506:LEU:HD13	1:B:515:LEU:HD12	1.67	0.77
1:A:507:LYS:HG2	1:A:507:LYS:O	1.84	0.76
1:B:428:VAL:HG12	1:B:439:ILE:HD11	1.67	0.75
1:B:323:GLN:O	1:B:326:LYS:HG3	1.86	0.75
1:A:88:ASN:O	1:A:92:THR:HG23	1.87	0.73
1:A:728:LEU:HD23	1:A:798:VAL:HG13	1.72	0.72
1:A:506:LEU:HD13	1:A:515:LEU:HD12	1.70	0.72
1:B:533:ARG:HE	1:B:554:ALA:H	1.36	0.71
1:B:76:GLU:CD	1:B:76:GLU:H	1.96	0.71
1:A:160:GLU:HG2	1:A:161:GLN:N	2.07	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:707:GLU:O	1:B:711:LYS:HG2	1.93	0.68
1:B:426:VAL:HG12	1:B:428:VAL:HG13	1.77	0.67
4:A:887:TDK:OM4	1:B:106:HIS:HE1	1.78	0.67
1:A:309:LYS:HG2	1:A:343:GLU:HG3	1.76	0.67
1:A:494:THR:OG1	1:A:692:LEU:HG	1.94	0.67
1:A:274:ILE:HG13	1:A:319:ASP:OD2	1.96	0.65
1:A:728:LEU:HD23	1:A:798:VAL:CG1	2.26	0.65
1:A:261:LEU:HD23	1:A:274:ILE:HD11	1.79	0.65
1:B:64:PRO:HB2	1:B:66:GLU:OE2	1.97	0.65
1:A:140:GLN:O	1:A:143:ILE:HG13	1.98	0.64
1:A:400:ALA:HB1	1:A:406:ALA:HA	1.78	0.64
1:A:434:GLU:H	1:A:434:GLU:CD	2.04	0.64
1:A:638:LEU:HD22	1:A:828:PHE:HB3	1.79	0.64
1:A:290:ILE:HD11	1:A:375:LYS:HD2	1.80	0.64
1:A:150:ARG:NH2	5:A:956:HOH:O	2.32	0.63
1:A:832:ASP:OD2	1:B:169:VAL:HB	1.98	0.63
1:A:859:ARG:HD3	1:A:861:GLU:OE2	1.99	0.63
1:A:344:THR:HA	1:A:347:LEU:HD12	1.81	0.62
1:A:301:LEU:HD13	1:A:351:TRP:CZ2	2.34	0.62
1:B:526:PHE:HB2	1:B:528:MET:HE2	1.80	0.62
1:A:177:TYR:CD2	1:A:192:VAL:HG11	2.35	0.62
1:A:300:GLU:HG3	1:A:301:LEU:N	2.15	0.61
1:A:261:LEU:HD12	1:A:323:GLN:HG2	1.82	0.61
1:A:536:GLY:HA2	1:A:557:LYS:HD2	1.82	0.60
1:A:297:ARG:HB3	1:A:359:LEU:HD23	1.83	0.60
1:A:414:MET:HE1	1:A:436:LEU:HD21	1.83	0.60
1:B:128:ARG:HG3	1:B:128:ARG:HH11	1.67	0.60
1:A:864:LYS:HE2	1:B:779:PRO:O	2.02	0.59
1:B:311:ILE:HA	1:B:314:MET:CE	2.32	0.59
1:B:326:LYS:HD3	1:B:391:ILE:HG23	1.84	0.59
1:A:304:LYS:HE2	1:A:347:LEU:HD23	1.84	0.59
1:B:287:TRP:CE3	1:B:385:VAL:HG22	2.37	0.58
1:A:309:LYS:HD2	1:A:341:TYR:CG	2.37	0.58
1:B:857:ALA:HB1	1:B:864:LYS:HD3	1.85	0.58
1:A:466:GLN:O	1:A:466:GLN:HG3	2.03	0.58
1:B:233:MET:O	1:B:239:LYS:HE3	2.04	0.58
1:A:821:ARG:HD2	1:A:855:GLU:CG	2.34	0.58
1:B:86:ARG:NH1	3:B:889:PO4:O2	2.37	0.58
1:B:863:ASP:OD1	1:B:865:LYS:HG3	2.05	0.57
1:A:536:GLY:HA3	1:A:557:LYS:NZ	2.18	0.57
1:B:323:GLN:HA	1:B:326:LYS:HE3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:MET:HE2	1:A:410:LYS:HA	1.86	0.57
1:A:502:LEU:HA	1:A:505:MET:CE	2.34	0.57
1:A:400:ALA:HB2	1:A:409:VAL:HG21	1.85	0.57
1:A:851:ALA:O	1:A:855:GLU:HG3	2.05	0.57
1:A:550:ARG:HH12	1:A:555:TYR:HB2	1.70	0.57
1:B:638:LEU:HD22	1:B:828:PHE:HB3	1.87	0.57
1:A:863:ASP:OD1	1:A:863:ASP:N	2.35	0.57
1:B:90:ILE:CG2	1:B:94:LEU:HD22	2.36	0.56
1:A:106:HIS:HE1	4:B:887:TDK:OM4	1.87	0.56
1:A:532:PHE:O	1:A:536:GLY:HA2	2.06	0.56
1:B:506:LEU:CD1	1:B:515:LEU:HD12	2.36	0.56
1:A:502:LEU:HA	1:A:505:MET:HE3	1.88	0.55
1:A:550:ARG:NH1	1:A:555:TYR:HB2	2.21	0.55
1:A:692:LEU:HD22	1:A:733:SER:HB3	1.88	0.55
1:A:863:ASP:O	1:A:866:VAL:HB	2.07	0.55
1:A:197:GLY:N	1:A:198:PRO:HD2	2.21	0.55
1:B:856:LEU:HB3	1:B:862:ILE:HG13	1.89	0.55
1:B:72:PRO:HG3	1:B:370:TYR:CE1	2.42	0.55
1:B:264:LEU:C	1:B:264:LEU:HD12	2.32	0.55
1:B:326:LYS:HE2	1:B:361:ARG:NH1	2.22	0.55
1:A:512:LYS:HG3	1:A:513:ASP:N	2.21	0.55
1:A:656:ILE:HD11	1:A:685:VAL:HG21	1.89	0.55
1:A:855:GLU:O	1:A:859:ARG:HG3	2.06	0.55
1:A:508:ASN:O	1:A:512:LYS:HB3	2.07	0.54
1:A:113:SER:HB3	1:A:258:ASN:ND2	2.22	0.54
1:B:177:TYR:HB3	1:B:178:PRO:CD	2.37	0.54
1:B:765:ALA:O	1:B:769:GLN:HG3	2.07	0.54
1:A:450:TYR:O	1:A:454:GLN:HG2	2.07	0.54
1:A:821:ARG:HD2	1:A:855:GLU:HG2	1.89	0.54
1:B:477:SER:O	1:B:480:ASP:HB2	2.07	0.54
1:B:664:TYR:CG	1:B:701:MET:HB2	2.42	0.54
1:B:532:PHE:O	1:B:536:GLY:HA2	2.07	0.53
1:B:140:GLN:O	1:B:143:ILE:HG13	2.09	0.53
1:B:326:LYS:HG2	1:B:361:ARG:NH2	2.22	0.53
1:B:197:GLY:N	1:B:198:PRO:HD2	2.24	0.53
1:B:859:ARG:HG3	1:B:861:GLU:HB2	1.90	0.53
1:A:126:ARG:HH22	1:A:132:ASP:CG	2.15	0.53
1:A:528:MET:O	1:A:531:LEU:HB2	2.10	0.52
1:B:533:ARG:NH1	1:B:533:ARG:HB3	2.24	0.52
1:B:400:ALA:HB2	1:B:409:VAL:HG21	1.92	0.52
1:B:533:ARG:NE	1:B:554:ALA:H	2.05	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:547:PRO:O	1:A:550:ARG:HB2	2.10	0.51
1:A:82:ARG:O	1:A:86:ARG:HG3	2.11	0.51
1:A:160:GLU:HG3	1:A:172:ASN:ND2	2.26	0.51
1:A:304:LYS:HE2	1:A:347:LEU:CD2	2.40	0.51
1:A:334:ARG:NE	1:A:353:ASP:OD1	2.44	0.51
1:A:539:SER:O	1:A:560:GLU:HA	2.11	0.50
1:B:128:ARG:HG2	1:B:133:GLY:HA2	1.93	0.50
1:A:420:ILE:HG13	1:A:421:ARG:N	2.26	0.50
1:A:622:PHE:CE1	1:A:677:MET:HE1	2.46	0.50
1:B:326:LYS:HE2	1:B:361:ARG:HH12	1.75	0.50
1:A:713:ILE:HB	1:A:764:LEU:HD11	1.92	0.50
5:A:893:HOH:O	1:B:192:VAL:HG22	2.10	0.50
1:A:294:TRP:HB3	1:A:298:TRP:CD1	2.47	0.50
1:A:729:LEU:HD11	1:A:792:MET:HE3	1.94	0.50
1:A:773:ARG:HD3	5:A:954:HOH:O	2.11	0.50
1:B:287:TRP:HE3	1:B:385:VAL:HG22	1.76	0.50
1:B:512:LYS:HG3	1:B:513:ASP:N	2.27	0.50
1:A:550:ARG:HD3	1:A:550:ARG:C	2.36	0.50
1:B:549:ASP:HB2	1:B:552:GLN:HG2	1.94	0.50
1:A:103:LEU:O	1:A:166:ARG:HD3	2.12	0.49
1:A:524:ARG:HG3	1:B:265:ASP:OD1	2.11	0.49
1:B:264:LEU:HD12	1:B:264:LEU:O	2.12	0.49
1:A:585:TYR:O	1:A:589:ASN:HA	2.13	0.49
1:A:177:TYR:HB3	1:A:178:PRO:CD	2.42	0.49
1:B:546:THR:HG23	1:B:546:THR:O	2.12	0.49
1:A:177:TYR:CG	1:A:192:VAL:HG11	2.48	0.49
1:B:539:SER:O	1:B:560:GLU:HA	2.12	0.49
1:B:811:VAL:O	1:B:815:VAL:HG23	2.13	0.49
1:B:128:ARG:HH11	1:B:128:ARG:CG	2.26	0.49
1:A:309:LYS:HD2	1:A:341:TYR:CD2	2.48	0.48
1:A:547:PRO:HD2	1:A:550:ARG:HG2	1.95	0.48
1:B:82:ARG:HG2	1:B:86:ARG:NH1	2.29	0.48
1:B:458:LEU:O	1:B:459:HIS:HB2	2.13	0.48
1:A:111:GLN:NE2	1:A:394:TYR:O	2.46	0.48
1:B:508:ASN:O	1:B:512:LYS:HB3	2.13	0.48
1:A:274:ILE:HG13	1:A:274:ILE:H	1.46	0.48
1:A:857:ALA:HB1	1:A:864:LYS:HG2	1.95	0.48
1:B:279:GLU:O	1:B:283:GLU:HB2	2.14	0.48
1:B:553:VAL:O	1:B:554:ALA:HB3	2.14	0.48
1:B:199:ILE:HG12	1:B:237:GLU:HB2	1.95	0.48
1:B:547:PRO:O	1:B:550:ARG:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:490:LYS:HD3	1:B:492:ILE:CG2	2.43	0.47
1:B:550:ARG:HA	1:B:550:ARG:HD2	1.59	0.47
4:A:887:TDK:OM1	4:A:887:TDK:H7'1	2.14	0.47
1:A:304:LYS:NZ	1:A:347:LEU:O	2.44	0.47
1:A:638:LEU:C	1:A:638:LEU:HD12	2.39	0.47
1:A:780:LEU:HD23	1:B:864:LYS:HB2	1.95	0.47
1:B:63:ILE:HG23	1:B:67:GLU:CD	2.39	0.47
1:B:569:ILE:HD11	4:B:887:TDK:H62	1.97	0.47
1:B:274:ILE:HG13	1:B:319:ASP:OD2	2.15	0.47
1:B:164:ASN:HB3	1:B:173:GLY:HA2	1.97	0.47
1:A:421:ARG:CG	1:A:428:VAL:HG13	2.45	0.46
1:B:713:ILE:HB	1:B:764:LEU:HD11	1.96	0.46
1:A:206:LYS:HD2	1:A:248:GLU:HG3	1.96	0.46
1:A:312:GLN:O	1:A:316:GLU:HG2	2.15	0.46
1:A:654:ASN:O	1:A:685:VAL:HG22	2.15	0.46
1:B:428:VAL:CG1	1:B:439:ILE:HD11	2.40	0.46
1:B:547:PRO:HG2	1:B:550:ARG:HD3	1.96	0.46
1:B:842:GLU:HA	1:B:847:TYR:CE2	2.50	0.46
1:A:352:THR:HG23	1:A:355:GLN:OE1	2.16	0.46
1:A:569:ILE:HD12	1:B:231:GLY:C	2.40	0.46
1:B:533:ARG:HB3	1:B:533:ARG:CZ	2.46	0.46
1:A:328:LYS:HG2	1:A:332:TYR:CD2	2.50	0.46
1:A:493:SER:HB2	1:A:692:LEU:HD12	1.98	0.46
1:A:536:GLY:O	1:A:564:ILE:HG23	2.16	0.46
1:B:177:TYR:HB3	1:B:178:PRO:HD2	1.97	0.46
1:A:400:ALA:HB1	1:A:406:ALA:CA	2.45	0.46
1:B:178:PRO:HA	1:B:187:TRP:CG	2.51	0.46
1:A:502:LEU:HD12	1:A:505:MET:HE3	1.97	0.45
1:A:531:LEU:HD12	1:A:531:LEU:HA	1.80	0.45
1:B:327:SER:C	1:B:328:LYS:HD2	2.41	0.45
1:B:506:LEU:O	1:B:512:LYS:NZ	2.48	0.45
1:B:800:SER:OG	1:B:843:VAL:HG22	2.16	0.45
1:B:245:ALA:HA	1:B:250:LEU:HD12	1.97	0.45
1:B:834:ARG:HD2	3:B:890:PO4:O3	2.15	0.45
1:B:842:GLU:HA	1:B:847:TYR:CD2	2.50	0.45
1:A:158:THR:OG1	1:A:160:GLU:CD	2.59	0.45
1:B:90:ILE:HG22	1:B:94:LEU:HD22	1.98	0.45
1:A:466:GLN:O	1:A:466:GLN:CG	2.65	0.45
1:A:543:GLN:NE2	1:A:547:PRO:HG3	2.32	0.45
1:A:878:ALA:HA	1:B:777:LEU:HG	1.99	0.45
1:B:844:ASP:HB2	5:B:1085:HOH:O	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:ARG:HD2	1:B:521:ASP:CG	2.42	0.45
1:A:143:ILE:O	1:A:143:ILE:HD12	2.17	0.45
1:A:458:LEU:O	1:A:459:HIS:HB2	2.17	0.45
1:A:466:GLN:HE22	1:A:589:ASN:HB2	1.80	0.45
1:A:857:ALA:HA	1:A:862:ILE:HG13	1.98	0.45
1:B:723:LYS:HD3	1:B:749:TYR:O	2.17	0.45
1:A:471:GLU:OE2	1:A:591:PRO:HD2	2.17	0.45
1:B:250:LEU:C	1:B:252:ASN:H	2.25	0.45
1:B:361:ARG:HG2	1:B:389:HIS:ND1	2.32	0.45
1:B:861:GLU:C	1:B:862:ILE:HG12	2.42	0.44
1:A:194:MET:HE1	5:A:1089:HOH:O	2.18	0.44
1:B:193:SER:HB3	1:B:196:LEU:HD12	1.99	0.44
1:A:314:MET:HA	1:A:322:TYR:OH	2.17	0.43
1:A:838:ARG:HB3	1:A:844:ASP:HB3	2.00	0.43
1:B:723:LYS:HE3	1:B:862:ILE:HD13	1.99	0.43
1:B:85:ILE:CG1	1:B:153:LEU:HD23	2.49	0.43
1:B:106:HIS:HD2	5:B:1142:HOH:O	2.00	0.43
1:B:719:ILE:HD13	1:B:742:ALA:HB1	2.01	0.43
1:A:144:SER:N	1:A:145:PRO:CD	2.82	0.43
1:A:323:GLN:HE22	1:A:392:LYS:H	1.65	0.43
1:A:638:LEU:C	1:A:638:LEU:CD1	2.91	0.43
1:B:213:HIS:ND1	1:B:560:GLU:OE1	2.51	0.43
1:A:195:GLY:O	1:A:198:PRO:HG2	2.19	0.43
1:B:773:ARG:HD3	5:B:970:HOH:O	2.18	0.43
1:B:856:LEU:HB3	1:B:861:GLU:HB3	2.00	0.43
1:B:539:SER:HB2	1:B:558:GLU:HG2	2.01	0.43
1:A:414:MET:HE1	1:A:436:LEU:CD2	2.48	0.43
1:B:63:ILE:HG23	1:B:67:GLU:OE2	2.19	0.43
1:B:851:ALA:O	1:B:855:GLU:HG3	2.19	0.43
1:A:374:LYS:HA	1:A:374:LYS:HD3	1.73	0.42
1:A:323:GLN:NE2	1:A:392:LYS:H	2.17	0.42
1:A:453:ALA:O	1:A:457:LYS:HD3	2.19	0.42
1:B:79:ARG:NH2	1:B:424:PHE:HA	2.34	0.42
1:B:90:ILE:HG22	1:B:91:MET:HE2	2.00	0.42
1:B:505:MET:CE	1:B:670:MET:HE2	2.49	0.42
1:B:864:LYS:HD3	1:B:864:LYS:N	2.34	0.42
4:B:887:TDK:OM1	4:B:887:TDK:N4'	2.45	0.42
1:A:711:LYS:HD2	1:A:787:TYR:CD1	2.54	0.42
1:A:725:LYS:NZ	1:A:754:ASP:OD2	2.51	0.42
1:A:328:LYS:HD3	1:A:328:LYS:N	2.33	0.42
1:B:177:TYR:CD2	1:B:192:VAL:HG11	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:549:ASP:O	1:B:552:GLN:HG2	2.19	0.42
1:B:125:PHE:HB3	1:B:462:LEU:HD21	2.01	0.42
1:B:237:GLU:CD	1:B:237:GLU:H	2.26	0.42
1:B:144:SER:N	1:B:145:PRO:CD	2.83	0.42
1:B:597:ILE:HA	1:B:624:ILE:O	2.20	0.42
1:A:520:ALA:C	1:A:521:ASP:O	2.59	0.42
1:A:564:ILE:HG13	1:A:565:LEU:N	2.32	0.42
1:A:635:GLY:HA3	1:B:103:LEU:O	2.19	0.42
1:B:66:GLU:CD	1:B:66:GLU:H	2.28	0.42
1:B:522:GLU:OE2	4:B:887:TDK:H6 ⁷	2.20	0.42
1:A:320:GLY:HA3	1:B:545:TYR:CD2	2.54	0.41
1:A:323:GLN:HA	1:A:323:GLN:OE1	2.20	0.41
1:A:512:LYS:HG3	1:A:513:ASP:H	1.85	0.41
1:A:538:TYR:HB2	1:A:562:GLY:O	2.19	0.41
1:B:81:ILE:HA	1:B:441:PHE:HZ	1.85	0.41
1:B:856:LEU:O	1:B:859:ARG:HG2	2.20	0.41
1:A:245:ALA:HA	1:A:250:LEU:HD12	2.01	0.41
1:A:578:TRP:CD1	1:A:594:PRO:HB2	2.55	0.41
1:A:821:ARG:HH12	1:A:854:GLY:HA3	1.85	0.41
1:B:471:GLU:OE2	1:B:591:PRO:HD2	2.20	0.41
1:B:519:ILE:O	1:B:519:ILE:HG12	2.20	0.41
1:B:585:TYR:O	1:B:589:ASN:HA	2.20	0.41
1:B:711:LYS:HA	1:B:711:LYS:HD2	1.72	0.41
1:B:845:ALA:O	1:B:849:VAL:HG23	2.18	0.41
1:B:436:LEU:N	1:B:437:PRO:CD	2.83	0.41
1:B:777:LEU:HD12	1:B:777:LEU:HA	1.88	0.41
1:A:273:LYS:HG2	5:A:1079:HOH:O	2.20	0.41
1:A:487:GLU:HG3	1:A:488:GLN:N	2.35	0.41
1:B:262:GLN:HA	1:B:267:PRO:HA	2.03	0.41
1:B:408:GLN:O	1:B:410:LYS:HE2	2.19	0.41
1:B:856:LEU:CB	1:B:862:ILE:HG13	2.51	0.41
1:A:884:ARG:NH2	1:B:770:ASP:OD2	2.54	0.41
1:B:511:ILE:HD12	1:B:515:LEU:HD21	2.02	0.41
1:A:471:GLU:HG2	5:A:1116:HOH:O	2.20	0.41
1:B:91:MET:O	1:B:95:ARG:HG3	2.20	0.41
1:B:317:THR:HB	1:B:322:TYR:CE2	2.55	0.41
1:B:664:TYR:CD1	1:B:701:MET:HB2	2.56	0.41
1:A:92:THR:HG22	1:A:438:TYR:HE1	1.85	0.41
1:A:711:LYS:HD3	5:A:1067:HOH:O	2.20	0.41
1:B:189:PHE:HA	1:B:190:PRO:HD3	1.78	0.41
1:B:196:LEU:HA	1:B:199:ILE:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:405:ILE:H	1:B:405:ILE:HG13	1.73	0.41
1:A:87:TRP:CD2	1:A:426:VAL:HG11	2.56	0.41
1:A:87:TRP:CE3	1:A:420:ILE:HD11	2.56	0.41
1:A:550:ARG:HD3	1:A:550:ARG:O	2.20	0.41
1:B:128:ARG:CG	1:B:128:ARG:NH1	2.83	0.41
1:B:177:TYR:CG	1:B:192:VAL:HG11	2.55	0.41
1:B:348:VAL:HA	1:B:351:TRP:CD1	2.56	0.41
1:B:567:GLU:HG3	1:B:574:ALA:HA	2.03	0.41
1:B:716:LEU:HD13	1:B:739:ARG:CZ	2.51	0.41
1:B:164:ASN:CB	1:B:173:GLY:HA2	2.50	0.41
1:A:250:LEU:C	1:A:252:ASN:H	2.30	0.40
1:A:410:LYS:HE3	1:A:410:LYS:HB3	1.84	0.40
1:A:716:LEU:HD13	1:A:739:ARG:CZ	2.52	0.40
1:A:768:GLY:HA2	1:A:786:PRO:HB3	2.02	0.40
1:B:81:ILE:HA	1:B:441:PHE:CZ	2.57	0.40
1:B:214:ARG:HD3	5:B:1098:HOH:O	2.20	0.40
1:A:627:THR:HB	1:A:633:LEU:HD13	2.03	0.40
1:A:726:VAL:O	1:A:753:SER:HA	2.22	0.40
1:A:177:TYR:HB2	1:A:192:VAL:HG21	2.04	0.40
1:A:341:TYR:HB2	1:A:344:THR:OG1	2.21	0.40
1:B:90:ILE:HG23	1:B:94:LEU:HD22	2.02	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	829/886 (94%)	785 (95%)	40 (5%)	4 (0%)	24	22
1	B	829/886 (94%)	790 (95%)	38 (5%)	1 (0%)	48	50
All	All	1658/1772 (94%)	1575 (95%)	78 (5%)	5 (0%)	36	36

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	553	VAL
1	A	347	LEU
1	A	550	ARG
1	A	398	ASP
1	B	400	ALA

5.3.2 Protein sidechains ⓘ

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	691/735 (94%)	616 (89%)	75 (11%)	6	4
1	B	691/735 (94%)	624 (90%)	67 (10%)	8	5
All	All	1382/1470 (94%)	1240 (90%)	142 (10%)	7	4

All (142) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	VAL
1	A	66	GLU
1	A	90	ILE
1	A	97	SER
1	A	98	LYS
1	A	126	ARG
1	A	150	ARG
1	A	153	LEU
1	A	160	GLU
1	A	168	GLU
1	A	192	VAL
1	A	209	LYS
1	A	216	LEU
1	A	228	LEU
1	A	249	LYS
1	A	274	ILE
1	A	277	GLU
1	A	278	LEU

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Mol	Chain	Res	Type
1	A	297	ARG
1	A	300	GLU
1	A	302	LEU
1	A	304	LYS
1	A	306	THR
1	A	323	GLN
1	A	328	LYS
1	A	335	GLU
1	A	340	LYS
1	A	343	GLU
1	A	350	ASP
1	A	354	GLU
1	A	356	ILE
1	A	367	LYS
1	A	375	LYS
1	A	382	LYS
1	A	405	ILE
1	A	410	LYS
1	A	418	ARG
1	A	420	ILE
1	A	428	VAL
1	A	439	ILE
1	A	443	GLU
1	A	446	GLU
1	A	457	LYS
1	A	464	SER
1	A	472	LYS
1	A	486	GLU
1	A	487	GLU
1	A	506	LEU
1	A	507	LYS
1	A	512	LYS
1	A	531	LEU
1	A	533	ARG
1	A	546	THR
1	A	550	ARG
1	A	551	GLU
1	A	553	VAL
1	A	555	TYR
1	A	560	GLU
1	A	564	ILE
1	A	638	LEU

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Mol	Chain	Res	Type
1	A	680	GLU
1	A	681	LYS
1	A	685	VAL
1	A	692	LEU
1	A	703	GLU
1	A	739	ARG
1	A	785	VAL
1	A	798	VAL
1	A	812	ARG
1	A	818	ASP
1	A	835	GLU
1	A	862	ILE
1	A	863	ASP
1	A	865	LYS
1	A	880	LYS
1	B	56	ILE
1	B	60	ILE
1	B	65	VAL
1	B	67	GLU
1	B	68	GLN
1	B	76	GLU
1	B	94	LEU
1	B	98	LYS
1	B	128	ARG
1	B	153	LEU
1	B	168	GLU
1	B	169	VAL
1	B	192	VAL
1	B	199	ILE
1	B	237	GLU
1	B	239	LYS
1	B	244	ILE
1	B	264	LEU
1	B	274	ILE
1	B	277	GLU
1	B	278	LEU
1	B	297	ARG
1	B	303	ARG
1	B	306	THR
1	B	309	LYS
1	B	326	LYS
1	B	328	LYS

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Mol	Chain	Res	Type
1	B	334	ARG
1	B	378	GLU
1	B	385	VAL
1	B	398	ASP
1	B	403	LYS
1	B	410	LYS
1	B	411	LYS
1	B	414	MET
1	B	428	VAL
1	B	430	ASP
1	B	433	ILE
1	B	456	GLN
1	B	472	LYS
1	B	475	LEU
1	B	479	GLN
1	B	484	LEU
1	B	490	LYS
1	B	506	LEU
1	B	512	LYS
1	B	519	ILE
1	B	550	ARG
1	B	558	GLU
1	B	560	GLU
1	B	577	SER
1	B	610	LEU
1	B	623	LEU
1	B	638	LEU
1	B	692	LEU
1	B	711	LYS
1	B	718	THR
1	B	720	GLU
1	B	777	LEU
1	B	818	ASP
1	B	843	VAL
1	B	858	LYS
1	B	859	ARG
1	B	861	GLU
1	B	862	ILE
1	B	864	LYS
1	B	873	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	HIS
1	A	131	GLN
1	A	159	GLN
1	A	170	HIS
1	A	172	ASN
1	A	448	HIS
1	A	466	GLN
1	A	503	ASN
1	A	534	GLN
1	A	552	GLN
1	A	588	ASN
1	A	618	GLN
1	A	737	HIS
1	A	810	GLN
1	B	106	HIS
1	B	131	GLN
1	B	419	HIS
1	B	543	GLN
1	B	737	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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	$\# Z > 2$
4	TDK	B	887	2	31,35,35	1.65	5 (16%)	41,55,55	1.42	6 (14%)
3	PO4	A	889	-	4,4,4	0.89	0	6,6,6	0.90	0
3	PO4	A	890	-	4,4,4	0.93	0	6,6,6	0.87	0
3	PO4	B	889	-	4,4,4	1.15	0	6,6,6	0.70	0
3	PO4	B	890	-	4,4,4	1.20	0	6,6,6	0.81	0
3	PO4	B	891	-	4,4,4	1.05	0	6,6,6	0.88	0
4	TDK	A	887	2	31,35,35	1.64	6 (19%)	41,55,55	1.39	5 (12%)

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
4	TDK	B	887	2	-	3/28/35/35	0/2/2/2
4	TDK	A	887	2	-	3/28/35/35	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	887	TDK	PC-OM3	-4.73	1.47	1.56
4	A	887	TDK	PC-OM3	-4.66	1.47	1.56
4	B	887	TDK	PB-O1B	3.29	1.60	1.50
4	A	887	TDK	PB-O1B	3.26	1.60	1.50
4	A	887	TDK	C6'-C5'	-2.95	1.32	1.37
4	B	887	TDK	C6'-C5'	-2.72	1.32	1.37
4	A	887	TDK	PA-O3A	2.63	1.62	1.59
4	B	887	TDK	PC-OM4	-2.25	1.47	1.51
4	B	887	TDK	PA-O3A	2.19	1.61	1.59
4	A	887	TDK	PC-OM4	-2.04	1.47	1.51
4	A	887	TDK	PB-O3B	-2.02	1.47	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	887	TDK	OM4-PC-CMA	-4.81	108.59	113.45
4	A	887	TDK	OM4-PC-CMA	-4.44	108.97	113.45
4	B	887	TDK	O3B-PB-O2B	2.91	118.70	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	887	TDK	O3B-PB-O2B	2.88	118.60	107.80
4	A	887	TDK	C6'-N1'-C2'	2.71	120.52	116.07
4	B	887	TDK	C6'-N1'-C2'	2.56	120.28	116.07
4	B	887	TDK	OM2-PC-OM4	-2.44	109.34	114.62
4	A	887	TDK	OM2-PC-OM4	-2.43	109.35	114.62
4	B	887	TDK	OM3-PC-OM4	2.32	117.01	111.17
4	A	887	TDK	OM3-PC-OM4	2.23	116.78	111.17
4	B	887	TDK	S1-C2-N3	-2.19	110.10	112.24

There are no chirality outliers.

All (6) torsion outliers are listed below:

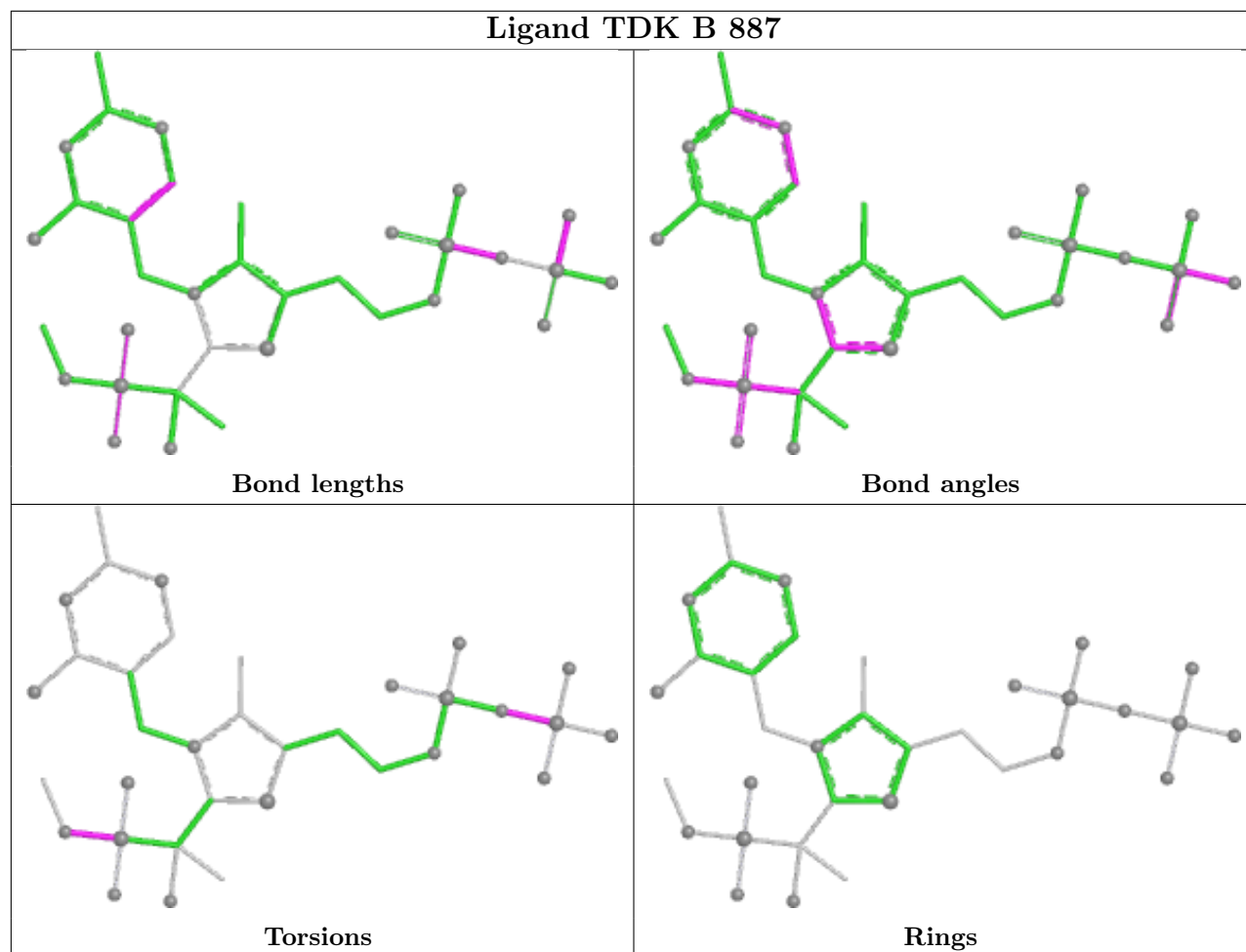
Mol	Chain	Res	Type	Atoms
4	A	887	TDK	PA-O3A-PB-O2B
4	B	887	TDK	PA-O3A-PB-O2B
4	A	887	TDK	CMC-OM2-PC-OM4
4	B	887	TDK	CMC-OM2-PC-OM4
4	A	887	TDK	CMC-OM2-PC-CMA
4	B	887	TDK	CMC-OM2-PC-CMA

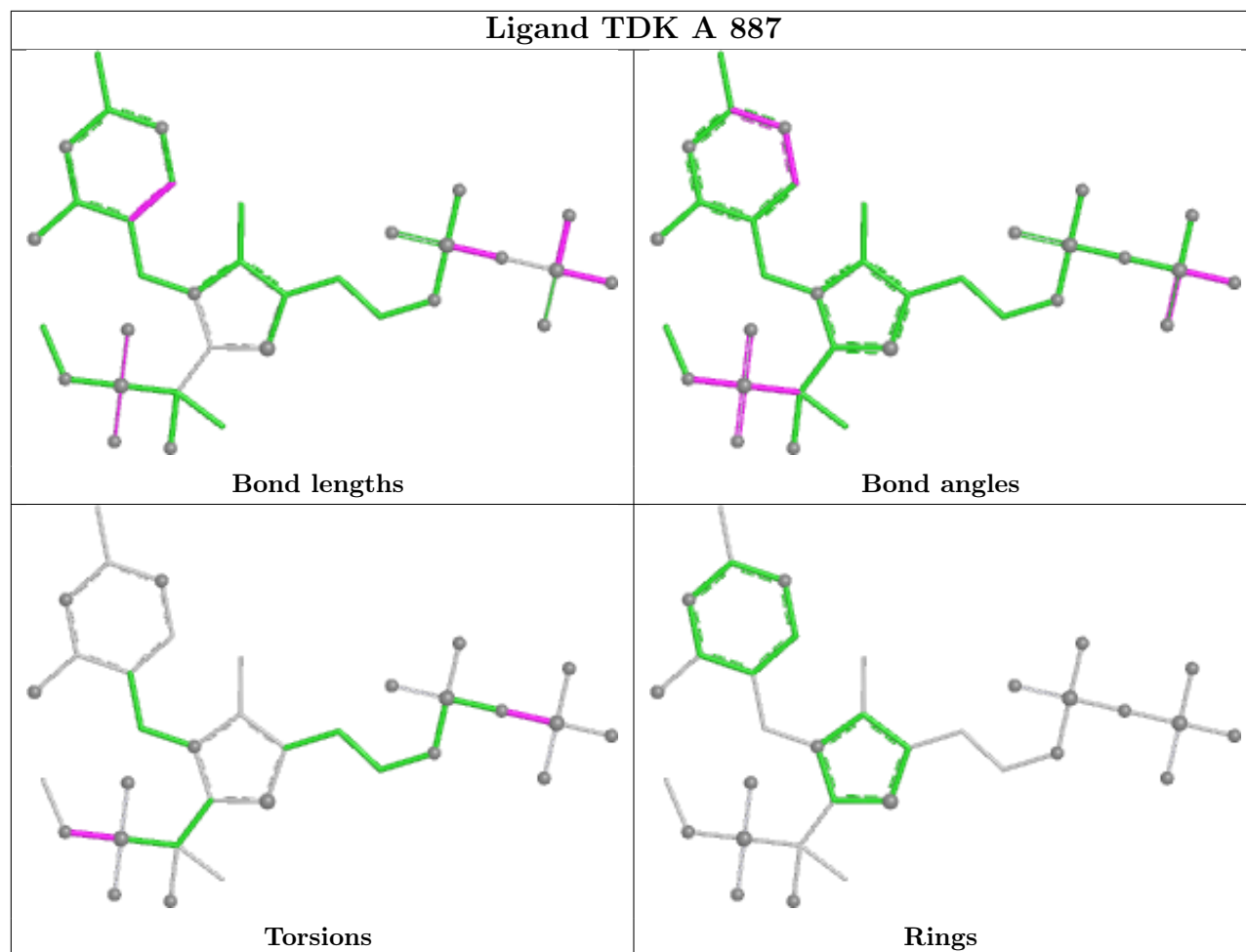
There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	887	TDK	4	0
3	B	889	PO4	1	0
3	B	890	PO4	1	0
4	A	887	TDK	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	831/886 (93%)	-0.48	10 (1%) 76 78	2, 8, 29, 43	0
1	B	831/886 (93%)	-0.60	3 (0%) 88 90	2, 6, 24, 37	0
All	All	1662/1772 (93%)	-0.54	13 (0%) 82 84	2, 7, 28, 43	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	818	ASP	3.5
1	A	549	ASP	3.2
1	A	552	GLN	3.1
1	A	541	ASN	3.0
1	A	398	ASP	2.8
1	A	818	ASP	2.5
1	A	323	GLN	2.4
1	A	352	THR	2.4
1	A	346	ALA	2.4
1	B	414	MET	2.3
1	B	67	GLU	2.3
1	A	555	TYR	2.1
1	A	551	GLU	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 oligosaccharides in this entry.

6.4 Ligands

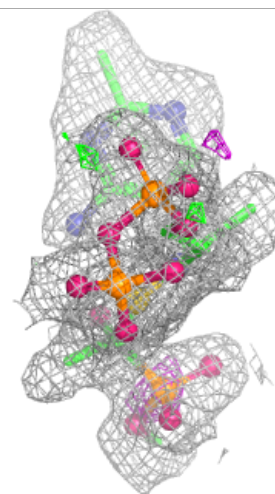
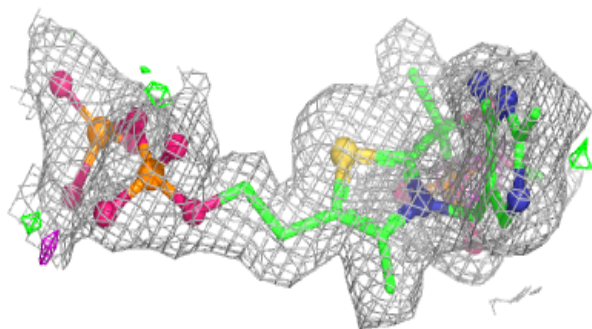
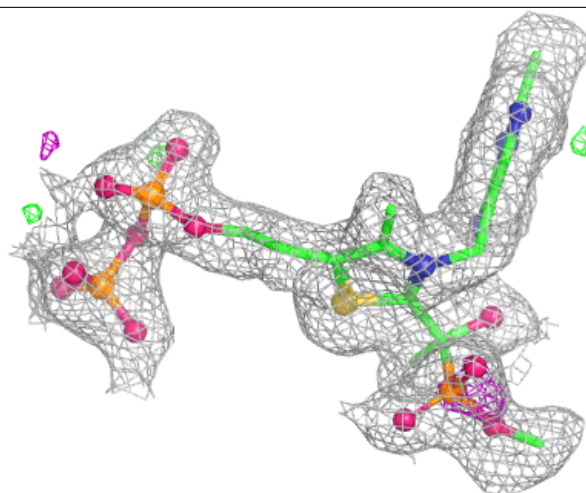
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	A	888	1/1	0.94	0.14	10,10,10,10	0
3	PO4	B	889	5/5	0.96	0.07	15,18,20,20	0
4	TDK	B	887	34/34	0.96	0.06	4,7,15,16	0
3	PO4	A	890	5/5	0.97	0.07	23,24,25,28	0
2	MG	B	888	1/1	0.97	0.15	13,13,13,13	0
3	PO4	B	890	5/5	0.97	0.07	23,24,25,27	0
3	PO4	B	891	5/5	0.97	0.07	24,24,27,28	0
4	TDK	A	887	34/34	0.97	0.05	2,3,13,14	0
3	PO4	A	889	5/5	0.97	0.05	12,13,14,16	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

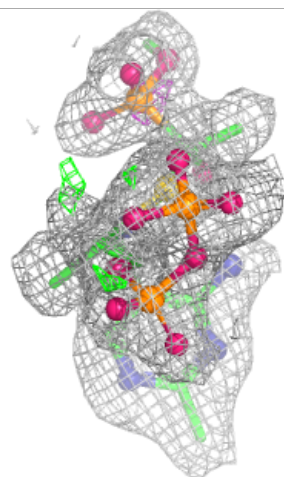
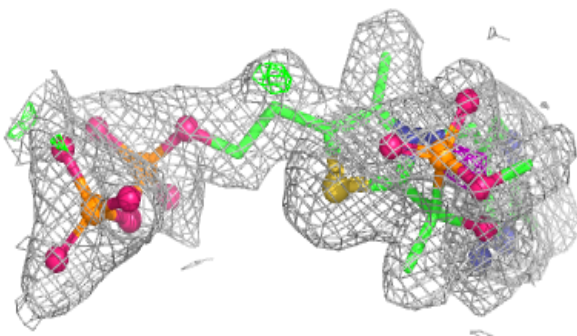
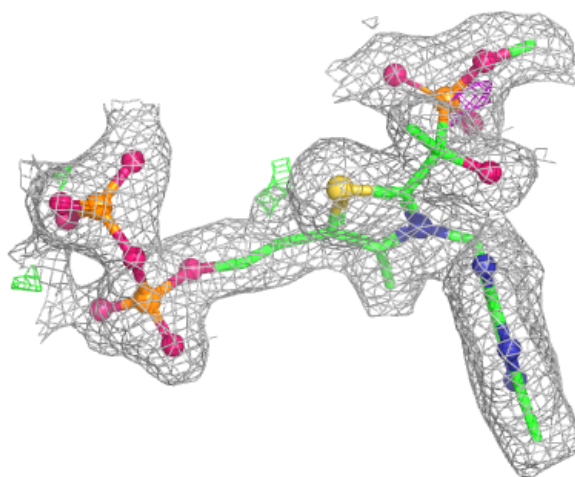
Electron density around TDK B 887:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around TDK A 887:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.