



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:58 AM UTC

PDB ID : 3LPL / pdb_00003lpl
Title : E. coli pyruvate dehydrogenase complex E1 component E571A mutant
Authors : Furey, W.
Deposited on : 2010-02-05
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

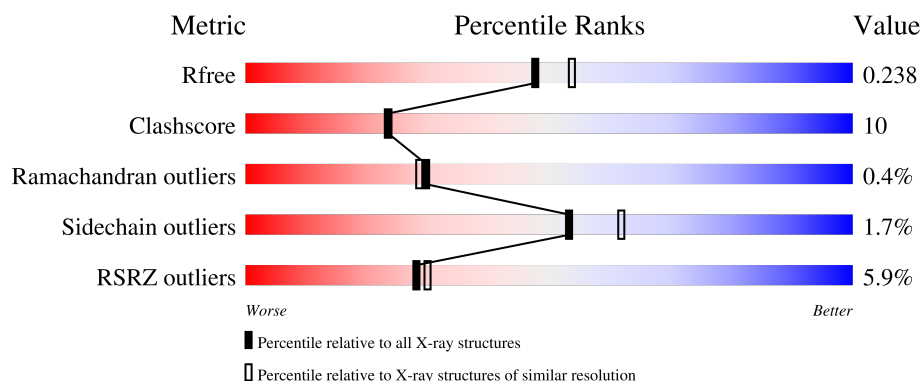
1 Overall quality at a glance

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	
1	B	886	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13497 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.

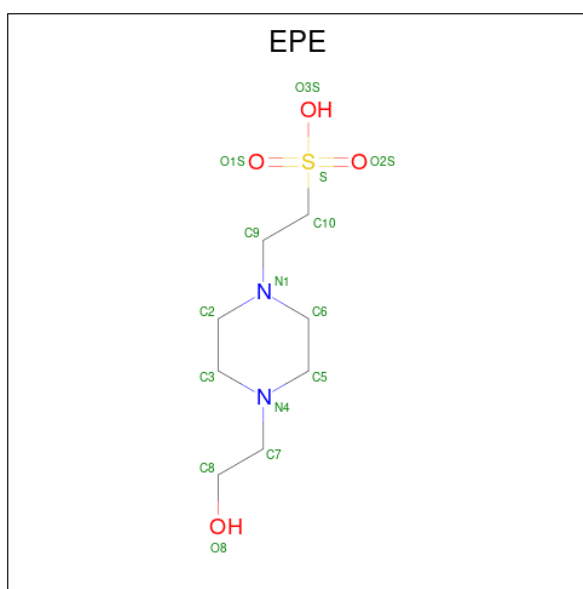
- Molecule 1 is a protein called Pyruvate dehydrogenase E1 component.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	801	Total	C	N	O	S	0	2	0
			6359	4031	1097	1205	26			
1	B	801	Total	C	N	O	S	0	1	0
			6349	4025	1094	1204	26			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	571	ALA	GLU	engineered mutation	UNP P0AFG9
B	571	ALA	GLU	engineered mutation	UNP P0AFG9

- Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: C₈H₁₈N₂O₄S).



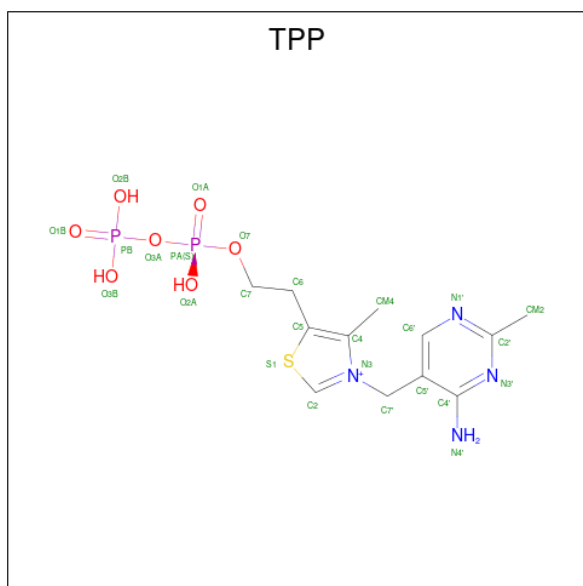
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

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

- Molecule 3 is THIAMINE DIPHOSPHATE (CCD ID: TPP) (formula: $C_{12}H_{19}N_4O_7P_2S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			11	2	7	2		
3	B	1	Total	C	O	P	0	0
			11	2	7	2		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



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

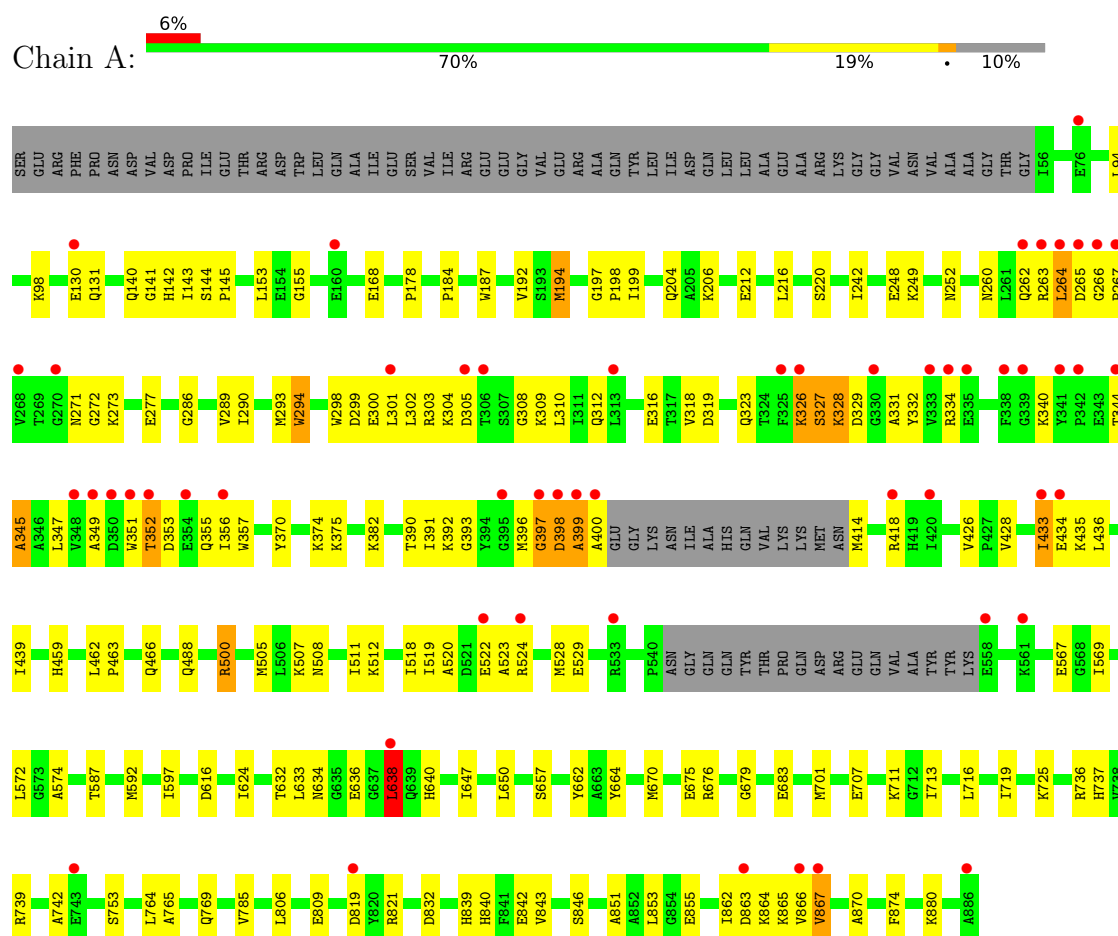
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	348	Total	O	0	0
			348	348		
6	B	377	Total	O	0	0
			377	377		

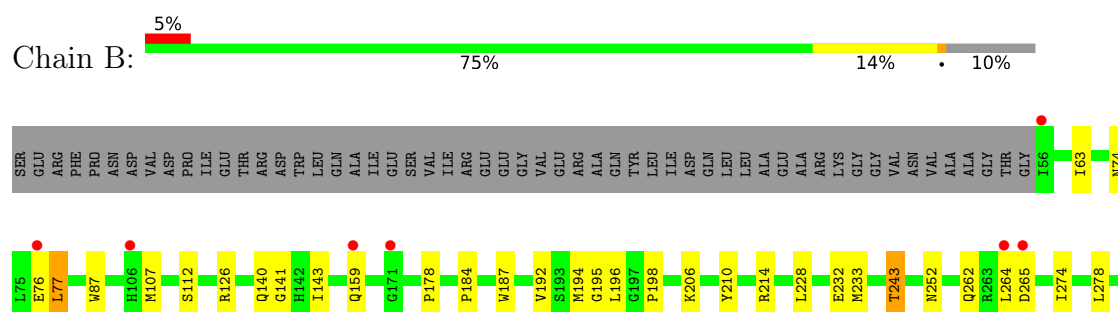
3 Residue-property plots [i](#)

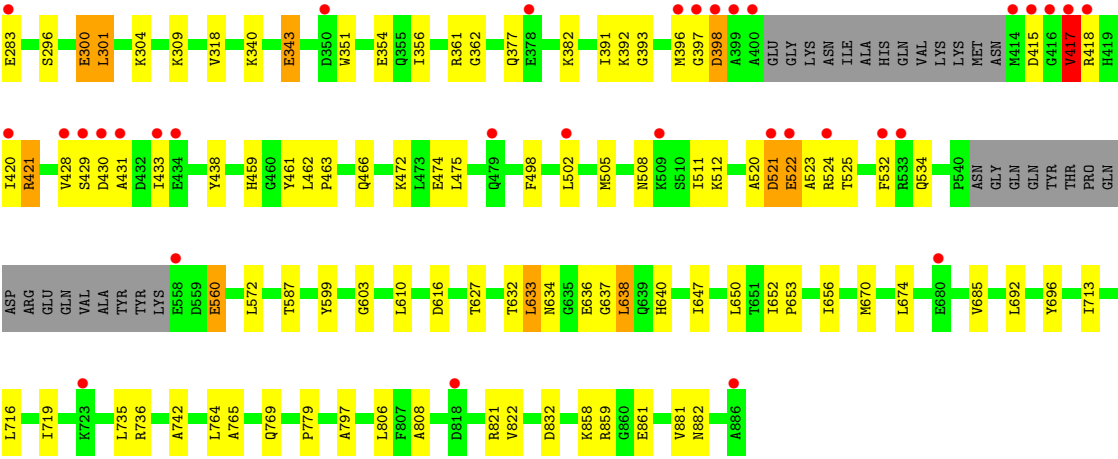
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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	81.82Å 142.22Å 82.52Å 90.00° 102.36° 90.00°	Depositor
Resolution (Å)	41.72 – 2.10 41.72 – 2.10	Depositor EDS
% Data completeness (in resolution range)	94.5 (41.72-2.10) 94.5 (41.72-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.72 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.197 , 0.238 0.203 , 0.238	Depositor DCC
R_{free} test set	5122 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.185	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13497	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% 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: TPP, MG, PO4, EPE

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	0.64	1/6504 (0.0%)	1.03	30/8794 (0.3%)
1	B	0.65	0/6493	1.02	21/8779 (0.2%)
All	All	0.65	1/12997 (0.0%)	1.03	51/17573 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	867	VAL	CA-CB	5.26	1.61	1.54

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	141	GLY	N-CA-C	8.01	122.33	112.64
1	B	63	ILE	N-CA-C	7.91	115.38	107.55
1	A	293	MET	N-CA-C	7.57	120.63	111.40
1	A	633	LEU	N-CA-C	-7.44	100.34	110.68
1	B	511	ILE	N-CA-C	7.40	120.00	112.90
1	B	534	GLN	N-CA-C	7.25	118.97	111.14
1	A	327	SER	N-CA-C	7.17	119.17	111.36
1	A	340	LYS	N-CA-C	-7.09	104.76	113.41
1	B	633	LEU	N-CA-C	-6.93	101.05	110.68
1	B	459	HIS	N-CA-C	6.76	121.48	113.23
1	A	459	HIS	N-CA-C	6.68	121.38	113.23
1	A	662	TYR	N-CA-C	6.67	120.47	110.14
1	A	141	GLY	N-CA-C	6.54	120.81	112.83
1	A	433	ILE	N-CA-C	6.53	118.06	110.62
1	B	636	GLU	N-CA-C	6.45	119.13	111.71
1	A	308	GLY	N-CA-C	-6.25	106.31	115.30
1	A	587	THR	N-CA-C	6.21	117.71	111.07
1	B	638	LEU	N-CA-C	6.20	121.60	111.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	636	GLU	N-CA-C	5.92	118.69	111.82
1	A	870	ALA	N-CA-C	-5.82	104.93	111.28
1	A	753	SER	N-CA-C	5.76	117.30	108.42
1	B	417	VAL	N-CA-C	-5.70	105.01	113.39
1	A	168	GLU	N-CA-C	5.64	120.32	113.38
1	B	637	GLY	N-CA-C	5.63	122.44	112.53
1	B	797	ALA	N-CA-C	5.62	118.71	109.72
1	B	461	TYR	N-CA-C	5.58	118.83	109.95
1	A	523	ALA	N-CA-C	-5.56	105.60	112.72
1	A	294	TRP	N-CA-C	5.55	118.28	109.50
1	A	242	ILE	N-CA-C	5.53	117.97	111.05
1	A	632	THR	N-CA-C	5.44	118.38	111.69
1	A	299	ASP	N-CA-C	5.42	117.94	111.71
1	B	252	ASN	N-CA-C	-5.40	105.91	112.88
1	B	356	ILE	N-CA-C	-5.38	105.29	110.72
1	B	283	GLU	N-CA-C	-5.33	105.37	111.07
1	B	296	SER	N-CA-C	5.30	117.80	111.71
1	A	511	ILE	N-CA-C	5.29	119.20	113.43
1	B	692	LEU	N-CA-C	5.28	118.19	109.85
1	B	632	THR	N-CA-C	5.26	118.16	111.69
1	B	393	GLY	N-CA-C	-5.25	105.83	114.76
1	A	289	VAL	N-CA-C	5.23	115.62	107.99
1	A	679	GLY	N-CA-C	-5.23	104.57	112.41
1	A	390	THR	N-CA-C	5.21	116.58	109.18
1	A	184	PRO	N-CA-C	5.20	120.69	113.98
1	A	843	VAL	CB-CA-C	-5.18	105.57	110.70
1	A	252	ASN	N-CA-C	-5.17	106.21	112.88
1	A	657	SER	N-CA-C	5.16	117.89	109.59
1	B	362	GLY	N-CA-C	5.11	119.06	112.83
1	A	345	ALA	N-CA-C	-5.09	106.32	112.54
1	B	184	PRO	N-CA-C	5.09	120.55	113.98
1	A	785	VAL	N-CA-C	5.05	113.49	107.73
1	A	204	GLN	N-CA-C	-5.04	105.86	111.36

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	6359	0	6192	146	0
1	B	6349	0	6186	127	0
2	A	15	0	17	0	0
2	B	15	0	18	0	0
3	A	11	0	2	0	0
3	B	11	0	2	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
6	A	348	0	0	3	0
6	B	377	0	0	13	0
All	All	13497	0	12417	256	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 10.

All (256) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:415:ASP:O	1:B:418:ARG:HG2	1.46	1.14
1:B:430:ASP:HA	1:B:433:ILE:HD12	1.47	0.96
1:A:519:ILE:HD11	1:A:528:MET:HE2	1.49	0.94
1:B:415:ASP:CG	1:B:418:ARG:HD3	1.93	0.94
1:B:421:ARG:HG3	1:B:421:ARG:HH11	1.32	0.94
1:A:863:ASP:O	1:A:866:VAL:HG12	1.68	0.93
1:B:430:ASP:HA	1:B:433:ILE:CD1	2.05	0.86
1:A:262:GLN:NE2	1:A:392:LYS:HD3	1.92	0.83
1:A:853:LEU:HB3	1:A:867:VAL:HG22	1.62	0.81
1:B:198:PRO:HG3	1:B:228:LEU:HD22	1.61	0.81
1:A:707:GLU:O	1:A:711:LYS:HG2	1.82	0.78
1:B:415:ASP:O	1:B:418:ARG:CG	2.32	0.77
1:A:863:ASP:OD2	1:A:865:LYS:HB2	1.85	0.76
1:A:414:MET:O	1:A:418:ARG:HG3	1.86	0.75
1:A:272:GLY:O	1:A:318:VAL:HG13	1.86	0.75
1:B:587:THR:HG22	6:B:911:HOH:O	1.86	0.74
1:A:806:LEU:HD11	1:B:806:LEU:HD11	1.70	0.74
1:A:140:GLN:O	1:A:143:ILE:HG13	1.87	0.73
1:A:263:ARG:O	1:A:392:LYS:NZ	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:736:ARG:NH2	1:A:737:HIS:NE2	2.36	0.72
1:A:301:LEU:HD21	1:A:351:TRP:CZ2	2.24	0.72
1:B:498:PHE:CZ	1:B:502:LEU:HD11	2.24	0.72
1:A:192:VAL:HG21	1:B:640:HIS:HE1	1.55	0.71
1:A:522:GLU:OE1	1:B:264:LEU:HD13	1.89	0.71
1:B:377:GLN:HG3	6:B:1085:HOH:O	1.91	0.71
1:B:76:GLU:CD	1:B:76:GLU:H	1.98	0.70
1:A:334:ARG:HB3	1:A:356:ILE:CD1	2.20	0.70
1:A:863:ASP:OD2	1:A:865:LYS:HE3	1.91	0.70
1:A:518:ILE:C	1:A:519:ILE:HD12	2.17	0.69
1:A:263:ARG:NH1	1:A:267:PRO:O	2.26	0.68
1:A:304:LYS:HE2	1:A:347:LEU:HD23	1.76	0.68
1:A:273:LYS:HE2	6:A:1156:HOH:O	1.94	0.67
1:A:277:GLU:OE2	1:B:243:THR:HG21	1.93	0.67
1:A:142[O]:HIS:ND1	1:A:142[O]:HIS:N	2.42	0.66
1:A:344:THR:HA	1:A:347:LEU:HD12	1.77	0.66
1:A:303:ARG:O	1:A:303:ARG:HG3	1.94	0.66
1:A:290:ILE:HD11	1:A:375:LYS:HD2	1.77	0.66
1:B:719:ILE:HD12	1:B:742:ALA:HB1	1.76	0.66
1:A:312:GLN:O	1:A:316:GLU:HG2	1.96	0.65
1:B:194:MET:HE1	3:B:887:TPP:H72	1.78	0.65
1:A:286:GLY:O	1:A:382:LYS:HE3	1.95	0.65
1:B:415:ASP:OD2	1:B:418:ARG:HD3	1.95	0.65
1:A:505:MET:HE1	1:A:670:MET:HE2	1.80	0.63
1:A:263:ARG:HG3	1:A:264:LEU:N	2.13	0.63
1:A:199:ILE:HD13	1:A:572:LEU:HG	1.80	0.63
1:A:853:LEU:CB	1:A:867:VAL:HG22	2.28	0.63
1:A:298:TRP:HA	1:A:301:LEU:HD12	1.80	0.63
1:B:421:ARG:HH11	1:B:421:ARG:CG	2.08	0.63
1:B:656:ILE:HD11	1:B:685:VAL:HG21	1.81	0.62
1:A:634:ASN:HB2	1:A:832:ASP:O	1.98	0.62
1:A:352:THR:OG1	1:A:355:GLN:HG3	2.00	0.62
1:B:300:GLU:HG3	1:B:301:LEU:N	2.16	0.61
1:B:194:MET:HE2	1:B:232:GLU:HB2	1.82	0.60
1:B:194:MET:HE3	6:B:1096:HOH:O	2.01	0.60
1:A:345:ALA:O	1:A:349:ALA:HB2	2.02	0.60
1:A:352:THR:HG23	1:A:355:GLN:OE1	2.01	0.60
1:A:426:VAL:CG1	1:A:439:ILE:HD11	2.31	0.60
1:B:192:VAL:HG22	1:B:192:VAL:O	2.01	0.60
1:B:522:GLU:HG3	1:B:599:TYR:HE1	1.66	0.60
1:B:421:ARG:C	1:B:421:ARG:HD2	2.27	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:LEU:HD23	1:A:310:LEU:HD23	1.83	0.59
1:A:433:ILE:HG22	1:A:434:GLU:OE2	2.01	0.59
1:A:192:VAL:HG22	1:A:192:VAL:O	2.03	0.59
1:A:840:HIS:HA	1:A:880:LYS:NZ	2.17	0.59
1:A:199:ILE:HD13	1:A:572:LEU:CD2	2.33	0.59
1:A:329:ASP:HA	1:A:357:TRP:CZ3	2.38	0.58
1:A:304:LYS:NZ	1:A:347:LEU:HA	2.17	0.58
1:B:696:TYR:HB3	1:B:736:ARG:HH11	1.67	0.58
1:B:421:ARG:HG3	1:B:421:ARG:NH1	2.09	0.58
1:A:524:ARG:HA	1:A:529:GLU:OE1	2.03	0.57
1:B:656:ILE:HG12	1:B:685:VAL:CG2	2.34	0.57
1:A:98:LYS:HD2	1:A:436:LEU:CD1	2.34	0.57
1:A:130:GLU:HB3	1:A:131:GLN:HE22	1.68	0.57
1:A:318:VAL:HG12	1:A:319:ASP:N	2.19	0.57
1:B:274:ILE:HG13	1:B:278:LEU:HG	1.87	0.57
1:B:415:ASP:C	1:B:417:VAL:H	2.13	0.57
1:B:421:ARG:HD2	1:B:421:ARG:O	2.04	0.57
1:B:159:GLN:HG3	1:B:438:TYR:CD2	2.40	0.56
1:B:194:MET:CE	1:B:232:GLU:HB2	2.35	0.56
1:A:397:GLY:O	1:A:398:ASP:C	2.48	0.56
1:A:505:MET:CE	1:A:670:MET:HE2	2.35	0.56
1:A:522:GLU:OE1	1:B:264:LEU:CD1	2.54	0.56
1:A:713:ILE:HB	1:A:764:LEU:HD11	1.88	0.56
1:B:318:VAL:HG22	6:B:1234:HOH:O	2.06	0.55
1:A:142[O]:HIS:HE1	1:A:194:MET:HE1	1.70	0.55
1:A:294:TRP:HB3	1:A:298:TRP:CD1	2.42	0.55
1:B:420:ILE:C	1:B:420:ILE:HD12	2.33	0.54
1:B:572:LEU:HD12	1:B:610:LEU:HD22	1.88	0.54
1:B:656:ILE:CG1	1:B:685:VAL:HG21	2.37	0.54
1:B:498:PHE:CE1	1:B:502:LEU:HD21	2.43	0.54
1:A:569:ILE:HD13	1:B:194:MET:HE1	1.90	0.54
1:A:638:LEU:HD22	1:B:616:ASP:CG	2.33	0.54
1:B:210:TYR:CZ	1:B:214:ARG:HD2	2.43	0.54
1:B:508:ASN:O	1:B:512:LYS:HB3	2.08	0.53
1:B:112:SER:HB3	1:B:392:LYS:HA	1.90	0.53
1:B:421:ARG:NH2	1:B:430:ASP:HB2	2.23	0.53
1:A:842:GLU:OE2	1:A:880:LYS:HE2	2.08	0.53
1:B:429:SER:C	1:B:431:ALA:N	2.67	0.53
1:B:472:LYS:HE3	1:B:474:GLU:OE2	2.09	0.53
1:B:140:GLN:O	1:B:143:ILE:HG13	2.09	0.53
1:A:98:LYS:HD2	1:A:436:LEU:HD12	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:819:ASP:OD2	1:A:855:GLU:HG2	2.09	0.52
1:B:498:PHE:CE2	1:B:502:LEU:HD11	2.44	0.52
1:B:421:ARG:CG	1:B:421:ARG:NH1	2.70	0.52
1:A:212:GLU:OE2	1:A:220:SER:HB3	2.10	0.52
1:A:507:LYS:HA	1:A:507:LYS:HE2	1.91	0.52
1:B:466:GLN:CD	6:B:952:HOH:O	2.52	0.52
1:B:107:MET:CE	1:B:396:MET:HE1	2.39	0.52
1:A:719:ILE:HD12	1:A:742:ALA:HB1	1.92	0.52
1:B:265:ASP:OD1	1:B:265:ASP:O	2.28	0.52
1:B:304:LYS:NZ	6:B:1242:HOH:O	2.40	0.52
1:B:429:SER:C	1:B:431:ALA:H	2.16	0.52
1:A:393:GLY:HA3	1:A:400:ALA:HB1	1.92	0.52
1:A:144:SER:OG	1:A:145:PRO:HD3	2.10	0.51
1:B:634:ASN:HB2	1:B:832:ASP:O	2.10	0.51
1:A:716:LEU:HD13	1:A:739:ARG:NH1	2.25	0.51
1:B:736:ARG:NH1	6:B:1233:HOH:O	2.33	0.51
1:A:206:LYS:HD2	1:A:248:GLU:HG3	1.92	0.51
1:A:262:GLN:HA	1:A:267:PRO:HA	1.92	0.51
1:B:107:MET:HE2	1:B:396:MET:HE1	1.93	0.51
1:B:206:LYS:NZ	6:B:1204:HOH:O	2.41	0.51
1:B:587:THR:HG21	6:B:970:HOH:O	2.10	0.51
1:A:318:VAL:CG1	1:A:319:ASP:N	2.74	0.51
1:A:396:MET:O	1:A:397:GLY:O	2.27	0.51
1:A:414:MET:HE2	1:A:434:GLU:HA	1.93	0.51
1:B:859:ARG:NH1	1:B:861:GLU:OE1	2.44	0.51
1:B:475:LEU:HD22	1:B:674:LEU:HG	1.93	0.51
1:B:198:PRO:HG3	1:B:228:LEU:CD2	2.35	0.51
1:B:656:ILE:CD1	1:B:685:VAL:HG21	2.40	0.51
1:A:508:ASN:O	1:A:512:LYS:HB3	2.11	0.50
1:A:192:VAL:CG2	1:B:640:HIS:HE1	2.23	0.50
1:A:396:MET:O	1:A:397:GLY:C	2.53	0.50
1:B:523:ALA:CB	1:B:532:PHE:HZ	2.25	0.50
1:B:206:LYS:HD2	6:B:1056:HOH:O	2.12	0.50
1:B:265:ASP:O	1:B:265:ASP:CG	2.54	0.50
1:B:159:GLN:NE2	1:B:438:TYR:HD2	2.10	0.50
1:A:334:ARG:HB3	1:A:356:ILE:HD13	1.94	0.49
1:B:126:ARG:HD3	6:B:1184:HOH:O	2.12	0.49
1:B:415:ASP:OD1	1:B:418:ARG:HD3	2.10	0.49
1:A:301:LEU:CD2	1:A:351:TRP:CZ2	2.96	0.49
1:A:519:ILE:CG2	1:A:520:ALA:N	2.76	0.49
1:A:370:TYR:OH	1:A:374:LYS:HE3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:ILE:HD12	1:B:143:ILE:C	2.37	0.49
1:B:808:ALA:HB3	1:B:822:VAL:CG1	2.42	0.49
1:A:716:LEU:HD13	1:A:739:ARG:CZ	2.42	0.49
1:A:199:ILE:HD13	1:A:572:LEU:CG	2.43	0.49
1:A:331:ALA:O	1:A:334:ARG:HG2	2.12	0.49
1:A:263:ARG:NE	1:B:521:ASP:OD1	2.45	0.49
1:A:616:ASP:OD2	1:B:638:LEU:HD22	2.13	0.48
1:B:195:GLY:C	1:B:198:PRO:HD2	2.38	0.48
1:B:656:ILE:HG12	1:B:685:VAL:HG21	1.95	0.48
1:A:326:LYS:HD2	1:A:391:ILE:HG23	1.95	0.48
1:B:398:ASP:OD1	1:B:398:ASP:N	2.46	0.48
1:B:505:MET:HE3	1:B:670:MET:HE2	1.95	0.48
1:A:192:VAL:HG21	1:B:640:HIS:CE1	2.43	0.48
1:B:178:PRO:HA	1:B:187:TRP:CG	2.49	0.48
1:B:262:GLN:NE2	1:B:392:LYS:HD3	2.29	0.48
1:A:353:ASP:O	1:A:356:ILE:N	2.47	0.47
1:A:142[O]:HIS:CE1	1:A:194:MET:HE1	2.49	0.47
1:B:524:ARG:HB3	1:B:524:ARG:CZ	2.44	0.47
1:A:262:GLN:CD	1:A:392:LYS:HD3	2.38	0.47
1:B:415:ASP:C	1:B:417:VAL:N	2.71	0.47
1:B:627:THR:HB	1:B:633:LEU:HD22	1.95	0.47
1:A:736:ARG:NH2	1:A:737:HIS:CE1	2.82	0.47
1:A:640:HIS:HE1	1:B:192:VAL:HG21	1.79	0.47
1:B:309:LYS:HG3	1:B:343:GLU:CG	2.44	0.47
1:B:524:ARG:HB3	1:B:524:ARG:NH1	2.29	0.47
1:A:332:TYR:O	1:A:332:TYR:CD1	2.68	0.46
1:A:664:TYR:CG	1:A:701:MET:HB2	2.49	0.46
1:B:417:VAL:CG2	1:B:433:ILE:CG2	2.93	0.46
1:B:656:ILE:HG12	1:B:685:VAL:HG22	1.97	0.46
1:A:178:PRO:HA	1:A:187:TRP:CG	2.50	0.46
1:A:304:LYS:CE	1:A:347:LEU:HD23	2.43	0.46
1:B:397:GLY:O	1:B:398:ASP:C	2.57	0.46
1:A:300:GLU:CD	1:A:303:ARG:NH1	2.73	0.46
1:A:862:ILE:HD12	1:A:866:VAL:CG1	2.45	0.46
1:A:334:ARG:NE	1:A:353:ASP:OD1	2.47	0.46
1:A:821:ARG:HH11	1:A:851:ALA:HA	1.81	0.46
1:A:434:GLU:H	1:A:434:GLU:CD	2.22	0.46
1:B:462:LEU:HB2	1:B:463:PRO:HA	1.98	0.45
1:B:808:ALA:HB3	1:B:822:VAL:HG11	1.98	0.45
1:B:821:ARG:NH2	1:B:858:LYS:HE2	2.31	0.45
1:A:263:ARG:CG	1:A:264:LEU:N	2.78	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:LEU:CD1	1:A:592:MET:HE1	2.46	0.45
1:A:846:SER:HB2	1:A:874:PHE:CG	2.52	0.45
1:B:765:ALA:O	1:B:769:GLN:HG3	2.16	0.45
1:A:131:GLN:NE2	1:A:131:GLN:N	2.64	0.45
1:A:216:LEU:HD11	1:A:592:MET:HE1	1.99	0.45
1:A:302:LEU:C	1:A:304:LYS:H	2.25	0.45
1:B:74:ASN:CG	1:B:77:LEU:HB2	2.42	0.45
1:B:192:VAL:O	1:B:192:VAL:CG2	2.65	0.44
1:A:290:ILE:CD1	1:A:375:LYS:HD2	2.46	0.44
1:A:466:GLN:NE2	6:A:1162:HOH:O	2.50	0.44
1:B:87:TRP:CZ3	1:B:421:ARG:HB3	2.52	0.44
1:A:304:LYS:HZ1	1:A:347:LEU:HA	1.82	0.44
1:A:326:LYS:O	1:A:357:TRP:HH2	2.00	0.44
1:B:233:MET:HB2	6:B:1150:HOH:O	2.17	0.44
1:B:713:ILE:HB	1:B:764:LEU:HD11	1.99	0.44
1:A:640:HIS:HE1	1:B:192:VAL:CG2	2.31	0.43
1:A:263:ARG:HG2	1:A:266:GLY:O	2.18	0.43
1:A:426:VAL:HG13	1:A:439:ILE:HD11	2.00	0.43
1:A:650:LEU:C	1:A:650:LEU:HD12	2.44	0.43
1:A:323:GLN:O	1:A:326:LYS:HB3	2.18	0.43
1:A:765:ALA:O	1:A:769:GLN:HG3	2.18	0.43
1:A:462:LEU:HB2	1:A:463:PRO:HA	2.01	0.43
1:B:429:SER:O	1:B:433:ILE:HG13	2.18	0.43
1:A:647:ILE:O	1:A:650:LEU:HG	2.19	0.43
1:A:153:LEU:C	1:A:155:GLY:H	2.26	0.43
1:B:523:ALA:CB	1:B:532:PHE:CZ	3.01	0.43
1:A:94:LEU:HD23	1:A:94:LEU:HA	1.87	0.43
1:B:301:LEU:HD13	1:B:351:TRP:CZ2	2.54	0.42
1:A:569:ILE:HD13	1:B:194:MET:CE	2.49	0.42
1:B:354:GLU:CD	1:B:354:GLU:H	2.27	0.42
1:A:300:GLU:OE1	1:A:303:ARG:NH1	2.52	0.42
1:B:87:TRP:CE3	1:B:420:ILE:HD11	2.54	0.42
1:B:361:ARG:HD2	1:B:391:ILE:HG13	2.01	0.42
1:A:426:VAL:HG12	1:A:428:VAL:HG23	2.02	0.42
1:A:676:ARG:HH21	1:A:683:GLU:CD	2.25	0.42
1:A:862:ILE:HD12	1:A:866:VAL:HG11	2.01	0.42
1:B:525:THR:O	1:B:525:THR:CG2	2.67	0.42
1:A:260:ASN:C	1:A:262:GLN:H	2.28	0.42
1:A:567:GLU:HG3	1:A:574:ALA:HA	2.01	0.42
1:A:597:ILE:HD13	1:A:624:ILE:CG2	2.50	0.42
1:B:309:LYS:HB2	1:B:343:GLU:HG3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:ASP:OD2	1:B:523:ALA:N	2.43	0.42
1:A:638:LEU:HD21	1:B:616:ASP:HB2	2.02	0.42
1:B:560:GLU:O	1:B:560:GLU:HG2	2.20	0.42
1:B:309:LYS:HG3	1:B:343:GLU:HG3	2.02	0.42
1:B:716:LEU:HD11	1:B:735:LEU:HD21	2.02	0.41
1:A:305:ASP:O	1:A:305:ASP:CG	2.61	0.41
1:A:326:LYS:O	1:A:357:TRP:CH2	2.74	0.41
1:A:839:HIS:HB3	6:A:1212:HOH:O	2.20	0.41
1:A:131:GLN:N	1:A:131:GLN:CD	2.78	0.41
1:A:327:SER:O	1:A:328:LYS:O	2.38	0.41
1:A:488:GLN:OE1	1:A:500:ARG:NH2	2.53	0.41
1:B:647:ILE:O	1:B:650:LEU:HG	2.20	0.41
1:B:195:GLY:O	1:B:198:PRO:HD2	2.20	0.41
1:A:300:GLU:OE2	1:A:303:ARG:NH1	2.54	0.41
1:B:498:PHE:CZ	1:B:502:LEU:HD21	2.55	0.41
1:A:326:LYS:HB3	1:A:326:LYS:NZ	2.36	0.41
1:B:523:ALA:HB1	1:B:532:PHE:CZ	2.56	0.41
1:B:652:ILE:HA	1:B:653:PRO:HD3	1.89	0.41
1:B:881:VAL:HG22	1:B:882:ASN:N	2.36	0.41
1:A:197:GLY:N	1:A:198:PRO:HD2	2.36	0.41
1:B:274:ILE:O	1:B:278:LEU:HG	2.20	0.41
1:A:304:LYS:HZ3	1:A:347:LEU:HA	1.86	0.40
1:A:248:GLU:O	1:A:249:LYS:C	2.65	0.40
1:B:520:ALA:O	1:B:522:GLU:N	2.54	0.40
1:B:603:GLY:HA3	6:B:1060:HOH:O	2.21	0.40
1:A:192:VAL:O	1:A:192:VAL:CG2	2.68	0.40
1:A:271:ASN:C	1:A:318:VAL:HG11	2.47	0.40
1:A:864:LYS:HD3	1:B:779:PRO:O	2.20	0.40
1:A:329:ASP:HA	1:A:357:TRP:CE3	2.57	0.40
1:A:397:GLY:O	1:A:399:ALA:N	2.54	0.40
1:A:806:LEU:HA	1:A:809:GLU:HB2	2.04	0.40
1:B:428:VAL:O	1:B:428:VAL:HG13	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	797/886 (90%)	761 (96%)	31 (4%)	5 (1%)	21	18
1	B	796/886 (90%)	755 (95%)	39 (5%)	2 (0%)	36	36
All	All	1593/1772 (90%)	1516 (95%)	70 (4%)	7 (0%)	30	28

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	328	LYS
1	A	398	ASP
1	A	399	ALA
1	A	397	GLY
1	B	521	ASP
1	B	522	GLU
1	A	638	LEU

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	666/734 (91%)	656 (98%)	10 (2%)	57	65
1	B	665/734 (91%)	653 (98%)	12 (2%)	51	60
All	All	1331/1468 (91%)	1309 (98%)	22 (2%)	53	62

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	194	MET
1	A	264	LEU
1	A	309	LYS
1	A	326	LYS
1	A	352	THR

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Mol	Chain	Res	Type
1	A	435	LYS
1	A	500	ARG
1	A	638	LEU
1	A	675	GLU
1	A	725	LYS
1	B	77	LEU
1	B	196	LEU
1	B	243	THR
1	B	300	GLU
1	B	301	LEU
1	B	340	LYS
1	B	343	GLU
1	B	382	LYS
1	B	398	ASP
1	B	417	VAL
1	B	421	ARG
1	B	560	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	GLN
1	A	159	GLN
1	A	172	ASN
1	A	262	GLN
1	A	456	GLN
1	A	466	GLN
1	A	534	GLN
1	A	588	ASN
1	A	695	ASN
1	B	419	HIS
1	B	468	ASN
1	B	697	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
5	PO4	B	890	-	4,4,4	0.56	0	6,6,6	0.78	0
5	PO4	A	890	-	4,4,4	0.46	0	6,6,6	0.65	0
3	TPP	A	887	4	9,10,27	1.74	2 (22%)	13,15,40	0.99	0
2	EPE	A	889	-	15,15,15	1.35	3 (20%)	19,20,20	2.15	6 (31%)
2	EPE	B	889	-	15,15,15	1.71	2 (13%)	19,20,20	2.11	5 (26%)
3	TPP	B	887	4	9,10,27	1.71	1 (11%)	13,15,40	0.95	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	EPE	A	889	-	-	3/9/19/19	0/1/1/1
2	EPE	B	889	-	-	0/9/19/19	0/1/1/1
3	TPP	A	887	4	-	3/10/10/17	-
3	TPP	B	887	4	-	4/10/10/17	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	889	EPE	O1S-S	4.18	1.56	1.45
3	B	887	TPP	PB-O1B	3.57	1.61	1.50
3	A	887	TPP	PB-O1B	3.51	1.61	1.50
2	B	889	EPE	C10-S	3.25	1.82	1.77
2	A	889	EPE	C10-S	3.02	1.81	1.77
2	A	889	EPE	O1S-S	2.83	1.53	1.45
2	A	889	EPE	C5-N4	2.30	1.53	1.46
3	A	887	TPP	PA-O3A	2.26	1.61	1.59

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	889	EPE	O3S-S-O2S	4.44	122.51	111.40
2	B	889	EPE	O3S-S-O2S	4.38	122.35	111.40
2	B	889	EPE	O2S-S-O1S	-3.88	101.20	113.82
2	A	889	EPE	O3S-S-O1S	-3.87	101.72	111.40
2	B	889	EPE	O3S-S-O1S	-3.75	102.02	111.40
2	A	889	EPE	O2S-S-O1S	-3.55	102.26	113.82
2	B	889	EPE	O2S-S-C10	3.37	111.82	106.73
2	A	889	EPE	O2S-S-C10	3.36	111.80	106.73
2	B	889	EPE	O3S-S-C10	3.34	112.54	106.00
2	A	889	EPE	C9-N1-C2	-3.09	103.01	111.24
2	A	889	EPE	C6-C5-N4	-2.28	106.05	110.65

There are no chirality outliers.

All (10) torsion outliers are listed below:

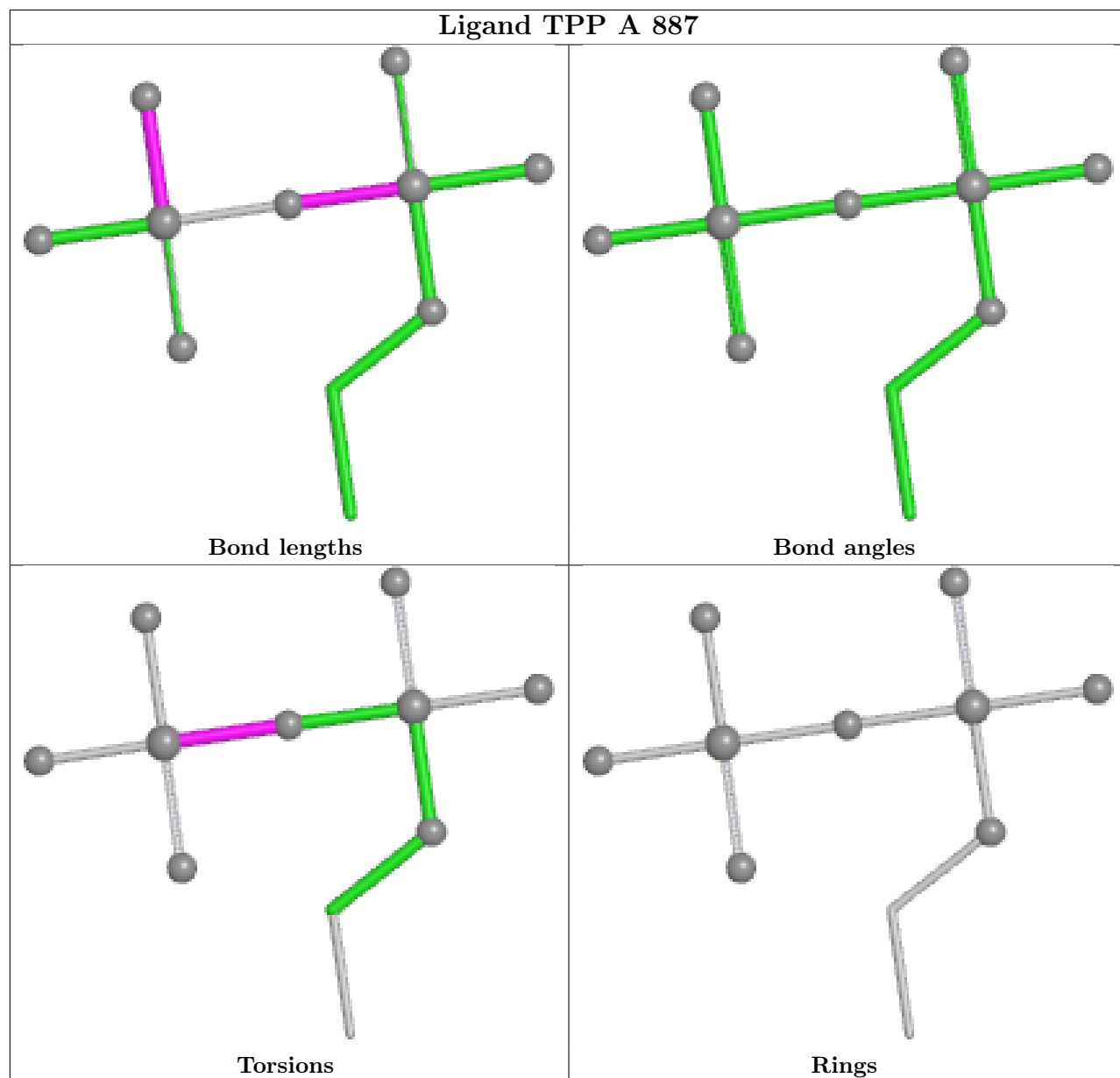
Mol	Chain	Res	Type	Atoms
2	A	889	EPE	C9-C10-S-O1S
2	A	889	EPE	C9-C10-S-O3S
3	A	887	TPP	PA-O3A-PB-O3B
2	A	889	EPE	C9-C10-S-O2S
3	A	887	TPP	PA-O3A-PB-O2B
3	B	887	TPP	C7-O7-PA-O1A
3	B	887	TPP	PA-O3A-PB-O1B
3	A	887	TPP	PA-O3A-PB-O1B
3	B	887	TPP	PA-O3A-PB-O2B
3	B	887	TPP	PA-O3A-PB-O3B

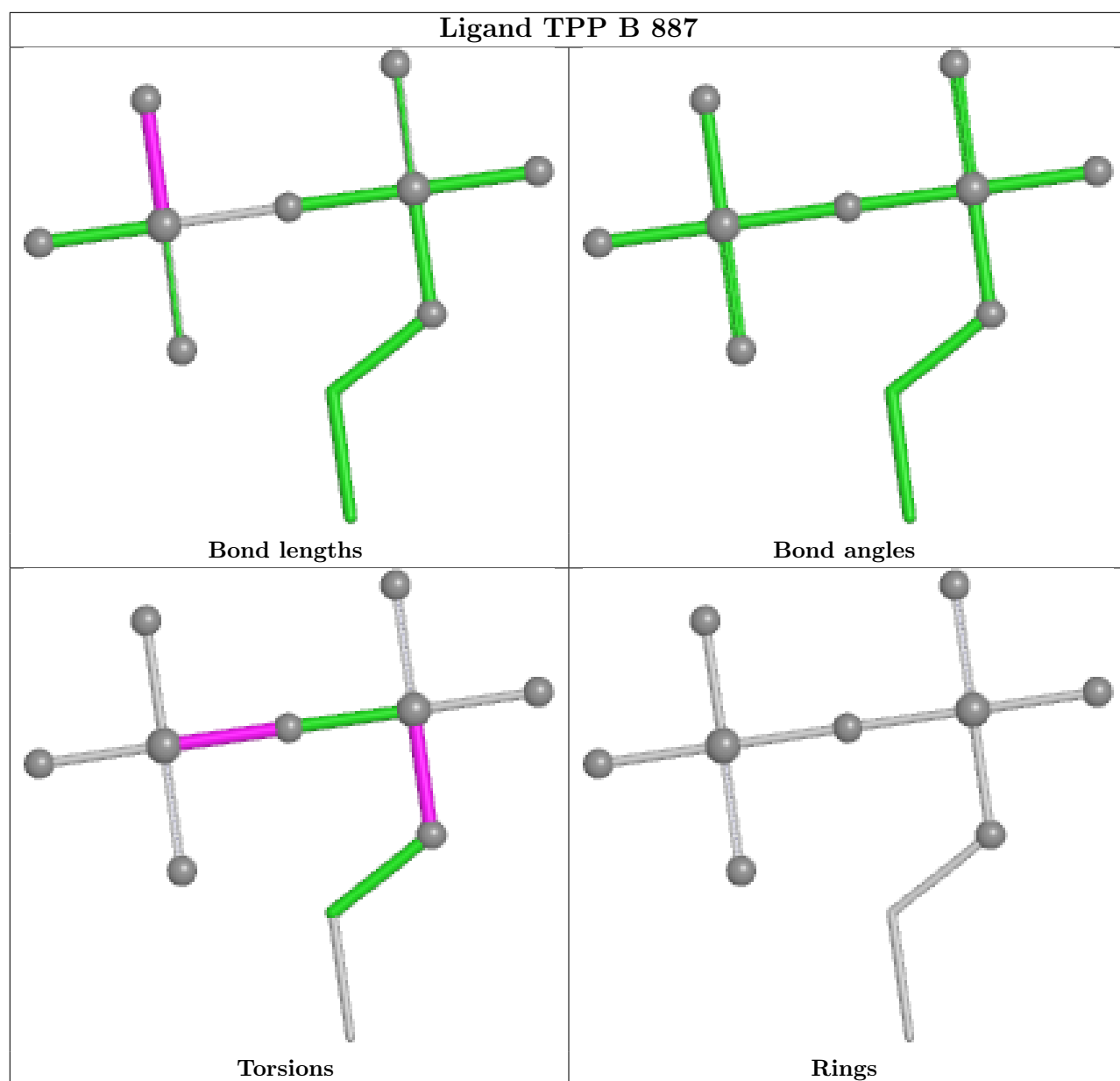
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	887	TPP	1	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

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%)	0.33	54 (6%) 24 25	11, 26, 40, 43	2 (0%)
1	B	801/886 (90%)	0.21	40 (4%) 34 36	11, 25, 35, 45	1 (0%)
All	All	1602/1772 (90%)	0.27	94 (5%) 28 30	11, 26, 38, 45	3 (0%)

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	399	ALA	9.3
1	B	400	ALA	7.6
1	B	416	GLY	6.7
1	B	414	MET	6.7
1	A	400	ALA	6.2
1	A	398	ASP	6.0
1	A	263	ARG	5.6
1	B	415	ASP	5.1
1	A	354	GLU	5.1
1	B	398	ASP	5.0
1	A	349	ALA	4.7
1	A	264	LEU	4.4
1	A	351	TRP	4.2
1	A	268	VAL	4.1
1	B	265	ASP	4.1
1	A	399	ALA	3.9
1	B	418	ARG	3.8
1	A	886	ALA	3.7
1	A	301	LEU	3.6
1	B	522	GLU	3.6
1	B	533	ARG	3.6
1	A	524	ARG	3.5
1	A	305	ASP	3.5
1	A	130	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	429	SER	3.5
1	B	818	ASP	3.4
1	A	350	ASP	3.3
1	B	417	VAL	3.3
1	B	521	ASP	3.3
1	A	333	VAL	3.2
1	B	434	GLU	3.2
1	B	428	VAL	3.2
1	A	265	ASP	3.1
1	B	350	ASP	3.1
1	A	341	TYR	3.1
1	A	867	VAL	3.1
1	A	266	GLY	3.0
1	B	397	GLY	3.0
1	A	160	GLU	3.0
1	B	283	GLU	3.0
1	A	397	GLY	3.0
1	B	171	GLY	3.0
1	B	378	GLU	3.0
1	B	264	LEU	3.0
1	B	502	LEU	2.9
1	A	418	ARG	2.9
1	A	342	PRO	2.9
1	A	326	LYS	2.9
1	A	76	GLU	2.8
1	A	819	ASP	2.8
1	B	524	ARG	2.7
1	B	723	LYS	2.7
1	B	532	PHE	2.7
1	B	433	ILE	2.7
1	B	420	ILE	2.6
1	B	431	ALA	2.6
1	A	325	PHE	2.6
1	A	638	LEU	2.5
1	A	330	GLY	2.5
1	A	262	GLN	2.4
1	A	743	GLU	2.4
1	A	267	PRO	2.4
1	A	335	GLU	2.4
1	A	866	VAL	2.4
1	B	396	MET	2.3
1	A	352	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	561	LYS	2.3
1	A	433	ILE	2.3
1	B	558	GLU	2.3
1	A	306	THR	2.3
1	B	886	ALA	2.3
1	B	76	GLU	2.3
1	B	159	GLN	2.3
1	B	479	GLN	2.3
1	B	56	ILE	2.2
1	A	863	ASP	2.2
1	B	430	ASP	2.2
1	A	434	GLU	2.2
1	A	338	PHE	2.2
1	A	348	VAL	2.2
1	A	339	GLY	2.2
1	B	680	GLU	2.1
1	A	395	GLY	2.1
1	B	106	HIS	2.1
1	A	313	LEU	2.1
1	A	522	GLU	2.1
1	A	356	ILE	2.1
1	A	420	ILE	2.1
1	A	533	ARG	2.1
1	B	509	LYS	2.1
1	A	334	ARG	2.1
1	A	558	GLU	2.0
1	A	344	THR	2.0
1	A	270	GLY	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 [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. The B-factors column lists the minimum,

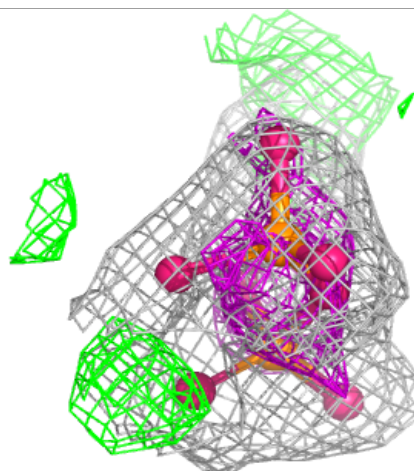
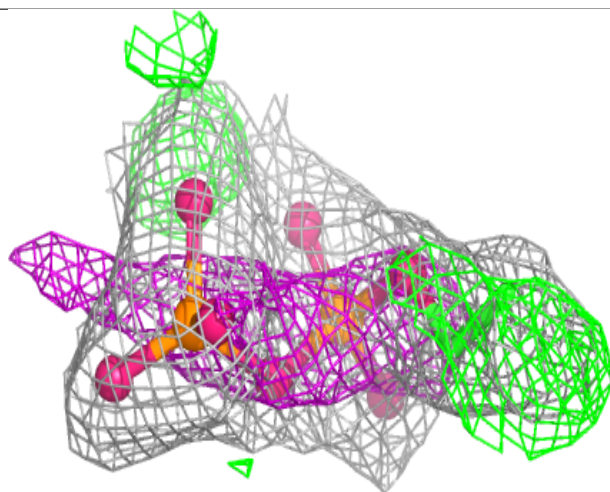
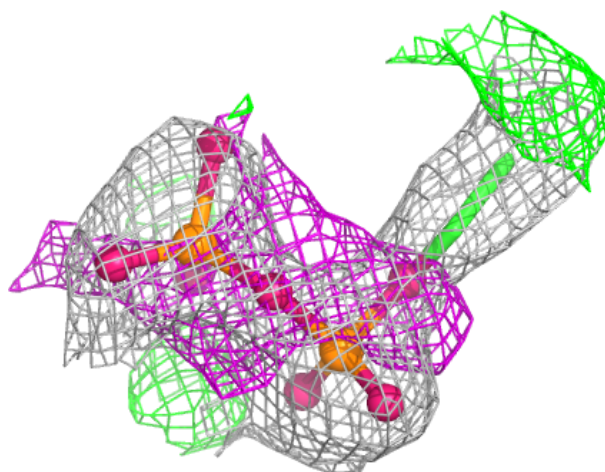
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	A	888	1/1	0.74	0.17	56,56,56,56	0
3	TPP	A	887	11/26	0.79	0.17	41,42,42,42	0
2	EPE	B	889	15/15	0.84	0.15	38,38,39,39	0
2	EPE	A	889	15/15	0.91	0.11	31,32,33,33	0
5	PO4	B	890	5/5	0.94	0.10	37,37,37,37	0
4	MG	B	888	1/1	0.95	0.05	32,32,32,32	0
5	PO4	A	890	5/5	0.95	0.10	36,36,36,37	0
3	TPP	B	887	11/26	0.95	0.08	35,35,35,35	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.

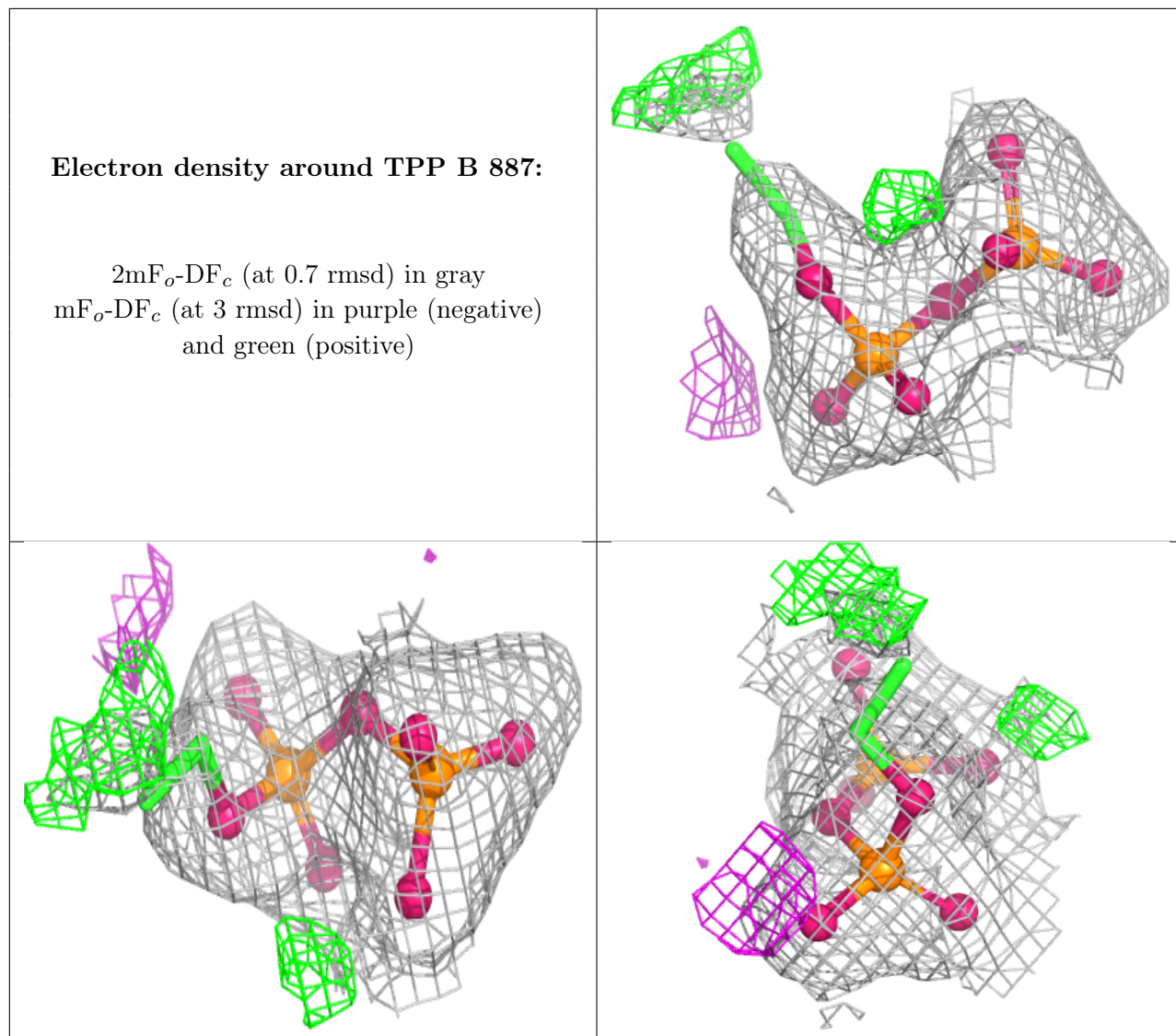
Electron density around TPP A 887:

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



Electron density around TPP 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.