



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:54 AM UTC

PDB ID : 8XHK / pdb_00008xhk
Title : Crystal structure of alpha-Oxoamine Synthase Alb29 with PLP cofactor
Authors : Xu, M.J.; Zhang, D.K.
Deposited on : 2023-12-18
Resolution : 2.79 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

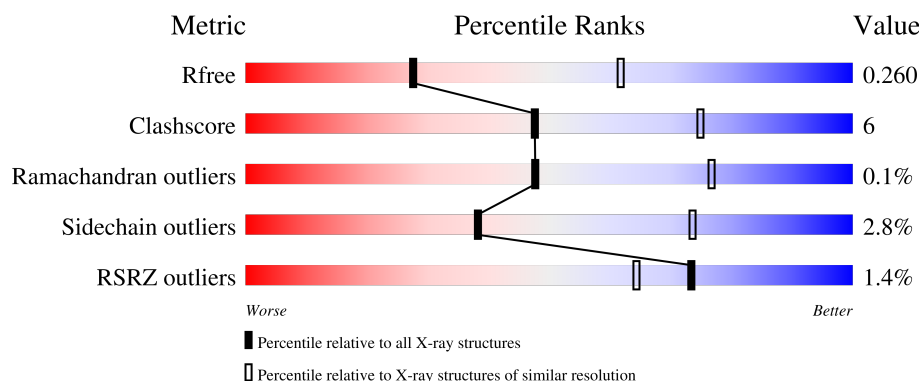
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	397	 86% 12% ..
1	B	397	 82% 16% ..
1	C	397	 86% 12% ..
1	D	397	 85% 13% ..
1	E	397	 84% 14% ..

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Mol	Chain	Length	Quality of chain
1	F	397	2% 81% 15% ••

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17952 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.

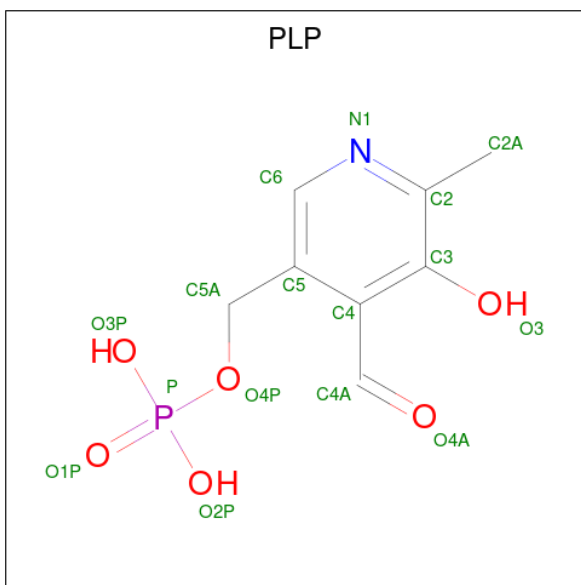
- Molecule 1 is a protein called Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	393	Total	C	N	O	S	0	0	0
			2975	1871	527	561	16			
1	A	394	Total	C	N	O	S	0	0	0
			2982	1877	529	560	16			
1	B	395	Total	C	N	O	S	0	0	0
			2991	1882	530	563	16			
1	D	395	Total	C	N	O	S	0	0	0
			2991	1882	530	563	16			
1	E	391	Total	C	N	O	S	0	0	0
			2958	1861	524	557	16			
1	F	390	Total	C	N	O	S	0	0	0
			2947	1853	520	558	16			

There are 6 discrepancies between the modelled and reference sequences:

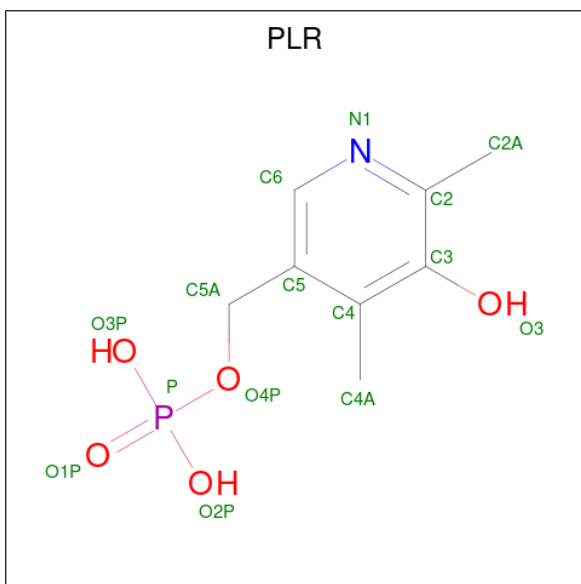
Chain	Residue	Modelled	Actual	Comment	Reference
C	87	GLY	SER	engineered mutation	UNP A0A6B9KSL0
A	87	GLY	SER	engineered mutation	UNP A0A6B9KSL0
B	87	GLY	SER	engineered mutation	UNP A0A6B9KSL0
D	87	GLY	SER	engineered mutation	UNP A0A6B9KSL0
E	87	GLY	SER	engineered mutation	UNP A0A6B9KSL0
F	87	GLY	SER	engineered mutation	UNP A0A6B9KSL0

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is (5-HYDROXY-4,6-DIMETHYLPYRIDIN-3-YL)METHYL DIHYDROGEN PHOSPHATE (CCD ID: PLR) (formula: $C_8H_{12}NO_5P$).



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

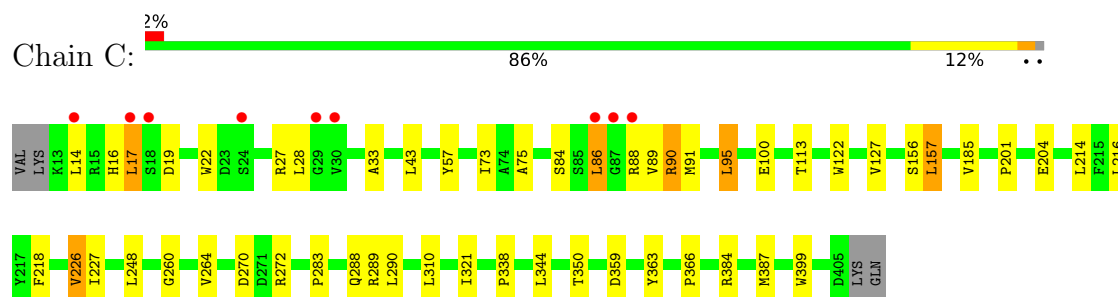
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	O	0	0
			1	1		
4	A	6	Total	O	0	0
			6	6		
4	B	5	Total	O	0	0
			5	5		
4	D	2	Total	O	0	0
			2	2		
4	E	1	Total	O	0	0
			1	1		
4	F	3	Total	O	0	0
			3	3		

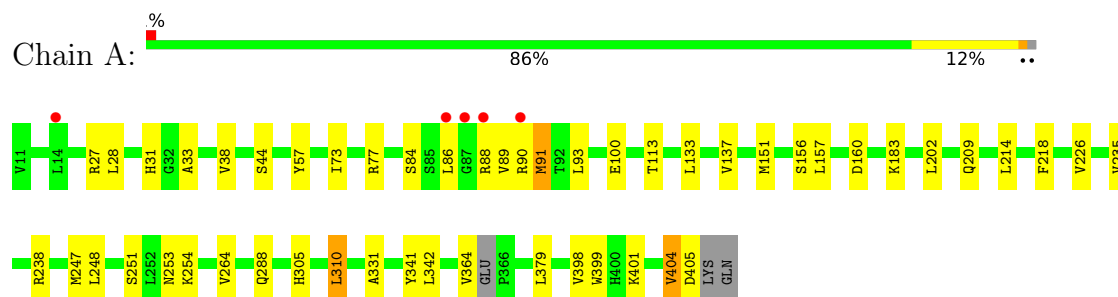
3 Residue-property plots [i](#)

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.

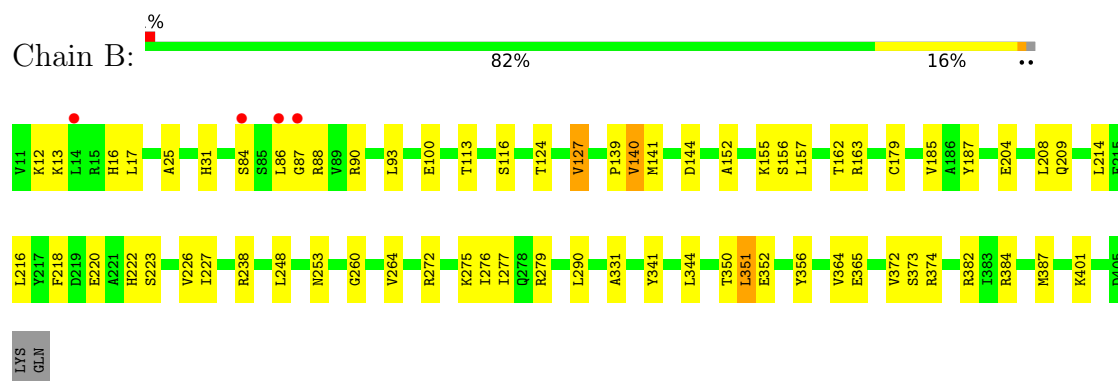
- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme



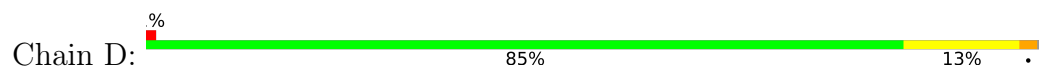
- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme

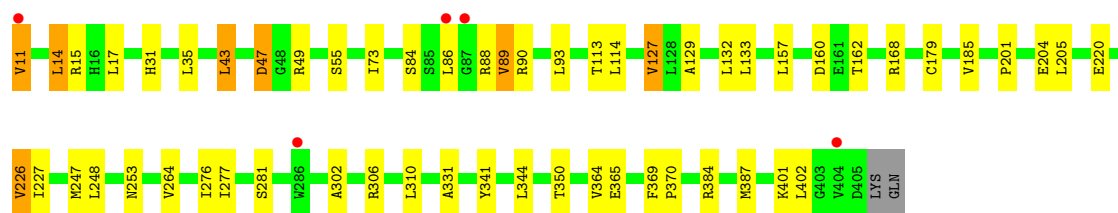


- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme

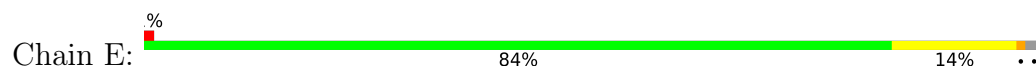


- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme

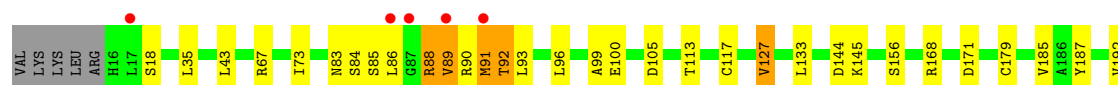
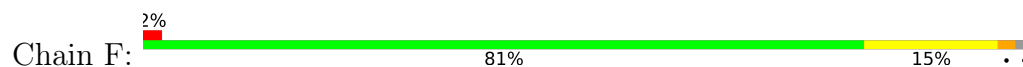




- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme



- Molecule 1: Aminotransferase class I/II-fold pyridoxal phosphate-dependent enzyme



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.11Å 110.59Å 354.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.19 – 2.79 25.19 – 2.79	Depositor EDS
% Data completeness (in resolution range)	95.7 (25.19-2.79) 95.5 (25.19-2.79)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.0.32, PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.203 , 0.255 0.213 , 0.260	Depositor DCC
R_{free} test set	3215 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	66.4	Xtriage
Anisotropy	0.566	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17952	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 28.65 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.7765e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PLR, PLP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.09	0/3036	0.25	0/4108
1	B	0.09	0/3046	0.26	0/4124
1	C	0.10	0/3030	0.27	0/4103
1	D	0.09	0/3046	0.27	0/4124
1	E	0.09	0/3013	0.25	0/4081
1	F	0.09	0/3002	0.27	0/4067
All	All	0.09	0/18173	0.26	0/24607

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	2982	0	2978	39	0
1	B	2991	0	2984	36	0
1	C	2975	0	2962	41	0
1	D	2991	0	2986	36	0
1	E	2958	0	2947	33	0
1	F	2947	0	2927	46	0
2	A	15	0	7	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	15	0	7	2	0
2	C	15	0	7	0	0
3	D	15	0	10	1	0
3	E	15	0	10	2	0
3	F	15	0	10	4	0
4	A	6	0	0	0	0
4	B	5	0	0	0	0
4	C	1	0	0	0	0
4	D	2	0	0	1	0
4	E	1	0	0	0	0
4	F	3	0	0	0	0
All	All	17952	0	17835	199	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 (199) 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:D:43:LEU:HD13	1:F:91:MET:HE1	1.63	0.80
1:F:226:VAL:HG23	1:F:227:ILE:HG13	1.67	0.77
1:C:86:LEU:HD12	1:A:253:ASN:HD22	1.55	0.71
1:A:84:SER:HA	1:A:93:LEU:HD12	1.72	0.70
1:C:75:ALA:HB1	1:C:95:LEU:HD11	1.76	0.68
1:B:384:ARG:HG3	1:B:387:MET:HG3	1.77	0.68
1:C:344:LEU:HD12	1:C:350:THR:HA	1.77	0.67
1:C:384:ARG:HG3	1:C:387:MET:HG3	1.76	0.67
1:E:89:VAL:HG23	1:E:91:MET:H	1.62	0.65
1:D:35:LEU:HD12	1:F:91:MET:HE3	1.80	0.64
1:B:84:SER:OG	1:B:290:LEU:O	2.15	0.62
1:D:84:SER:HA	1:D:93:LEU:HD12	1.81	0.62
1:F:86:LEU:HB2	1:F:91:MET:HG3	1.81	0.61
1:D:89:VAL:HG11	1:F:363:TYR:HE2	1.66	0.61
1:F:220:GLU:OE1	1:F:223:SER:OG	2.17	0.60
1:E:226:VAL:HG13	1:E:227:ILE:HG13	1.83	0.60
1:C:157:LEU:HD13	1:A:157:LEU:HD23	1.84	0.59
1:B:127:VAL:HG21	1:B:277:ILE:HG12	1.84	0.59
1:E:84:SER:HA	1:E:93:LEU:HD12	1.83	0.59
1:C:88:ARG:HA	1:A:57:TYR:CD1	2.38	0.58
1:B:226:VAL:HG13	1:B:227:ILE:HG13	1.84	0.58
1:D:160:ASP:OD2	1:F:156:SER:OG	2.20	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:344:LEU:HD12	1:D:350:THR:HA	1.85	0.58
1:B:179:CYS:HA	1:B:185:VAL:HG21	1.86	0.57
1:D:248:LEU:HB2	1:D:264:VAL:HB	1.86	0.57
1:F:84:SER:HA	1:F:93:LEU:HD12	1.87	0.56
1:C:43:LEU:HD22	1:A:86:LEU:HD22	1.87	0.56
1:C:363:TYR:HE2	1:A:89:VAL:HG11	1.70	0.56
1:D:14:LEU:HD12	1:D:15:ARG:HD2	1.88	0.56
1:A:86:LEU:HD12	1:A:91:MET:HB2	1.87	0.56
1:C:344:LEU:HD21	1:C:399:TRP:HH2	1.71	0.55
1:A:27:ARG:HD2	1:A:28:LEU:HD22	1.88	0.55
1:D:226:VAL:HG22	1:D:227:ILE:HG13	1.90	0.54
1:C:156:SER:OG	1:A:160:ASP:OD2	2.20	0.53
1:B:162:THR:OG1	1:B:163:ARG:N	2.40	0.53
1:A:137:VAL:HG11	1:A:183:LYS:HD2	1.89	0.53
1:C:84:SER:H	1:C:86:LEU:HD23	1.73	0.52
1:F:331:ALA:HB3	1:F:341:TYR:HB3	1.91	0.52
1:E:248:LEU:HB2	1:E:264:VAL:HB	1.92	0.52
1:C:57:TYR:CD2	1:A:88:ARG:HB3	2.45	0.52
1:E:133:LEU:HD22	1:E:273:LYS:HG2	1.92	0.51
1:D:88:ARG:HH22	3:F:501:PLR:P	2.34	0.51
1:C:33:ALA:H	1:A:91:MET:HE2	1.75	0.51
1:B:144:ASP:OD2	1:B:187:TYR:OH	2.28	0.50
1:D:31:HIS:CE1	1:F:90:ARG:HD3	2.46	0.50
1:F:127:VAL:HG21	1:F:277:ILE:HG12	1.93	0.50
1:A:133:LEU:HD13	1:A:247:MET:HE1	1.94	0.50
1:F:88:ARG:HH11	1:F:88:ARG:HB3	1.75	0.50
1:F:89:VAL:HB	1:F:91:MET:SD	2.52	0.50
1:B:365:GLU:OE2	1:B:382:ARG:NH2	2.45	0.49
1:A:100:GLU:HG2	1:A:113:THR:HG23	1.94	0.49
1:F:168:ARG:HG3	1:F:171:ASP:HB2	1.93	0.49
1:C:226:VAL:HG22	1:C:227:ILE:HG13	1.95	0.49
1:E:260:GLY:HA3	1:E:290:LEU:HD21	1.95	0.49
1:C:86:LEU:HB2	1:C:288:GLN:HB2	1.94	0.49
1:D:127:VAL:HG21	1:D:277:ILE:HG12	1.94	0.49
1:C:321:ILE:HD11	1:C:338:PRO:HB3	1.95	0.49
1:B:344:LEU:HD12	1:B:350:THR:HA	1.93	0.49
1:A:57:TYR:CD2	1:A:254:LYS:HE3	2.48	0.49
1:B:84:SER:HA	1:B:93:LEU:HD12	1.95	0.49
1:D:35:LEU:HD12	1:F:91:MET:CE	2.43	0.49
1:E:134:THR:HG21	1:E:139:PRO:HG3	1.94	0.49
1:F:192:VAL:HG22	1:F:199:LEU:HD22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:ASP:HA	1:C:22:TRP:HD1	1.77	0.49
1:E:134:THR:OG1	1:E:137:VAL:O	2.27	0.49
1:D:55:SER:HB2	1:F:89:VAL:HG11	1.93	0.48
1:C:366:PRO:O	1:A:90:ARG:NH2	2.46	0.48
1:A:251:SER:OG	1:A:253:ASN:OD1	2.32	0.48
1:D:89:VAL:HG12	1:D:90:ARG:H	1.79	0.48
1:E:216:LEU:HB3	1:E:218:PHE:CE2	2.48	0.48
1:A:88:ARG:NH2	1:A:288:GLN:HB3	2.29	0.48
1:B:209:GLN:HA	1:B:214:LEU:H	1.79	0.48
1:F:254:LYS:HZ1	3:F:501:PLR:C4	2.27	0.48
1:C:90:ARG:HD2	1:C:90:ARG:H	1.78	0.47
1:A:248:LEU:HB2	1:A:264:VAL:HB	1.96	0.47
1:E:344:LEU:HD12	1:E:350:THR:HA	1.95	0.47
1:F:224:THR:HG1	1:F:250:THR:HG1	1.50	0.47
1:F:272:ARG:O	1:F:276:ILE:HG12	2.14	0.47
1:D:47:ASP:OD2	1:D:49:ARG:NE	2.43	0.47
1:F:133:LEU:HD13	1:F:247:MET:HE1	1.96	0.47
1:D:73:ILE:HG23	1:F:73:ILE:HG23	1.96	0.47
1:D:86:LEU:HD22	1:F:43:LEU:HD22	1.95	0.47
1:B:124:THR:HG22	1:B:277:ILE:HD13	1.96	0.47
1:E:344:LEU:HD21	1:E:399:TRP:HH2	1.78	0.47
1:B:331:ALA:HB3	1:B:341:TYR:HB3	1.96	0.47
1:F:344:LEU:HD12	1:F:350:THR:HA	1.96	0.47
1:A:331:ALA:HB3	1:A:341:TYR:HB3	1.97	0.47
1:C:84:SER:H	1:C:86:LEU:CD2	2.27	0.47
1:C:90:ARG:HG3	1:A:31:HIS:CE1	2.50	0.47
1:B:248:LEU:HB2	1:B:264:VAL:HB	1.97	0.47
1:D:302:ALA:O	1:D:306:ARG:HG3	2.14	0.47
1:A:202:LEU:HB2	1:A:238:ARG:HH21	1.79	0.46
1:F:85:SER:HB2	1:F:88:ARG:HA	1.97	0.46
1:D:11:VAL:HA	1:D:14:LEU:HD11	1.96	0.46
1:C:248:LEU:HB2	1:C:264:VAL:HB	1.97	0.46
1:E:254:LYS:HZ2	3:E:501:PLR:C4	2.28	0.46
1:C:270:ASP:OD2	1:C:272:ARG:NH1	2.49	0.46
1:B:204:GLU:N	1:B:204:GLU:OE1	2.48	0.46
1:F:100:GLU:HG2	1:F:113:THR:HG23	1.98	0.46
1:C:216:LEU:HB3	1:C:218:PHE:CE2	2.51	0.46
1:E:321:ILE:HD11	1:E:338:PRO:HB3	1.98	0.46
1:A:305:HIS:HA	1:A:310:LEU:HD13	1.98	0.46
1:E:116:SER:HB2	3:E:501:PLR:O2P	2.16	0.46
1:B:351:LEU:HD13	1:B:351:LEU:HA	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:129:ALA:HB1	1:D:162:THR:HG22	1.97	0.46
1:E:127:VAL:HG21	1:E:277:ILE:HG12	1.97	0.45
1:A:398:VAL:HA	1:A:401:LYS:HE2	1.99	0.45
1:C:73:ILE:HG23	1:A:73:ILE:HG23	1.99	0.45
1:C:86:LEU:HD12	1:A:253:ASN:ND2	2.25	0.45
1:C:19:ASP:HA	1:C:22:TRP:CD1	2.52	0.45
1:B:222:HIS:NE2	2:B:501:PLP:O3	2.50	0.45
1:B:87:GLY:HA3	1:E:56:SER:HA	1.98	0.45
1:F:179:CYS:HA	1:F:185:VAL:HG21	1.99	0.45
1:C:73:ILE:HD13	1:A:77:ARG:HG2	1.99	0.45
1:C:260:GLY:HA3	1:C:290:LEU:HD21	1.99	0.45
1:B:152:ALA:O	1:B:155:LYS:HG3	2.16	0.45
1:A:209:GLN:HA	1:A:214:LEU:H	1.83	0.44
1:F:202:LEU:HB2	1:F:238:ARG:HH21	1.82	0.44
1:C:88:ARG:HB3	1:C:89:VAL:H	1.62	0.44
1:B:100:GLU:HG2	1:B:113:THR:HG23	1.99	0.44
1:B:88:ARG:HB2	1:E:57:TYR:CD1	2.52	0.44
1:E:129:ALA:HB1	1:E:162:THR:HG22	1.99	0.44
1:F:204:GLU:N	1:F:204:GLU:OE1	2.50	0.44
1:A:404:VAL:HG13	1:A:405:ASP:H	1.82	0.44
1:E:209:GLN:HA	1:E:214:LEU:H	1.82	0.44
1:D:133:LEU:HD13	1:D:247:MET:HE1	1.98	0.44
1:D:201:PRO:O	1:D:205:LEU:HG	2.17	0.44
1:B:25:ALA:HB1	1:B:31:HIS:HB2	1.99	0.44
1:E:316:LYS:HE3	1:E:320:ASN:HD21	1.81	0.44
1:E:190:ASP:OD1	1:E:190:ASP:N	2.51	0.44
1:E:369:PHE:CD1	1:E:370:PRO:HA	2.53	0.44
1:C:122:TRP:HB3	1:C:283:PRO:HG2	1.99	0.44
1:F:260:GLY:HA3	1:F:290:LEU:HD21	2.00	0.44
1:D:132:LEU:HB3	1:D:276:ILE:HG21	1.99	0.43
1:D:365:GLU:HB3	1:F:89:VAL:HG12	2.00	0.43
1:E:128:LEU:HG	1:E:141:MET:HE1	2.00	0.43
1:C:204:GLU:OE1	1:C:204:GLU:N	2.44	0.43
1:E:86:LEU:HD12	1:E:91:MET:HB3	2.01	0.43
1:F:305:HIS:HA	1:F:310:LEU:HD13	2.00	0.43
1:C:86:LEU:HD22	1:C:289:ARG:O	2.18	0.43
1:D:331:ALA:HB3	1:D:341:TYR:HB3	2.01	0.43
1:C:157:LEU:HD12	1:A:156:SER:HB2	2.00	0.43
1:B:140:VAL:HG13	1:B:185:VAL:HG13	2.00	0.43
1:E:92:THR:HB	1:E:96:LEU:HD23	2.01	0.43
1:B:139:PRO:HG2	1:B:141:MET:HE3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:253:ASN:ND2	1:E:88:ARG:HD2	2.33	0.43
1:D:113:THR:C	1:D:114:LEU:HD12	2.44	0.43
1:C:100:GLU:HG2	1:C:113:THR:HG23	2.00	0.43
1:F:144:ASP:OD2	1:F:187:TYR:OH	2.36	0.43
1:A:364:VAL:HG11	1:A:379:LEU:HB3	2.00	0.42
1:D:384:ARG:HG3	1:D:387:MET:HG3	2.00	0.42
1:F:92:THR:HG23	1:F:96:LEU:HD23	2.00	0.42
1:F:99:ALA:HA	1:F:299:ILE:HG12	2.01	0.42
1:E:172:VAL:HG11	1:E:204:GLU:HB3	2.01	0.42
1:B:12:LYS:HG3	1:B:13:LYS:H	1.83	0.42
1:A:218:PHE:CG	1:A:235:VAL:HG21	2.54	0.42
1:A:398:VAL:HG13	1:A:401:LYS:HE2	2.00	0.42
1:B:372:VAL:HG22	1:B:373:SER:H	1.85	0.42
1:D:204:GLU:OE1	1:D:204:GLU:N	2.46	0.42
1:D:253:ASN:HD21	1:F:88:ARG:HE	1.67	0.42
1:F:18:SER:HB2	1:F:370:PRO:HD3	2.01	0.42
1:F:117:CYS:HB3	3:F:501:PLR:O4P	2.19	0.42
1:A:89:VAL:HG12	1:A:90:ARG:H	1.85	0.42
1:B:216:LEU:HB3	1:B:218:PHE:CE2	2.55	0.42
1:C:359:ASP:O	1:B:374:ARG:NH1	2.53	0.42
1:A:342:LEU:HD23	1:A:399:TRP:CZ2	2.55	0.42
1:B:156:SER:OG	1:E:160:ASP:OD2	2.28	0.41
1:B:260:GLY:HA3	1:B:290:LEU:HD21	2.02	0.41
1:C:185:VAL:O	1:C:214:LEU:HD12	2.21	0.41
1:D:253:ASN:HB2	1:F:83:ASN:OD1	2.20	0.41
1:D:369:PHE:CD1	1:D:370:PRO:HA	2.55	0.41
3:F:501:PLR:H4A3	3:F:501:PLR:H5A1	1.89	0.41
1:A:89:VAL:HG12	1:A:90:ARG:N	2.35	0.41
1:D:17:LEU:HD23	1:D:17:LEU:HA	1.95	0.41
1:E:56:SER:HB3	1:E:384:ARG:HD3	2.02	0.41
1:C:91:MET:HE1	1:A:33:ALA:HB3	2.02	0.41
1:B:275:LYS:O	1:B:279:ARG:HG2	2.20	0.41
1:F:384:ARG:HG3	1:F:387:MET:HG3	2.02	0.41
1:F:105:ASP:OD1	1:F:268:ARG:NH2	2.49	0.41
1:B:116:SER:HB2	2:B:501:PLP:O1P	2.20	0.41
1:E:100:GLU:HG2	1:E:113:THR:HG23	2.03	0.41
1:F:335:ASN:HB3	1:F:337:VAL:H	1.85	0.41
1:C:16:HIS:CG	1:C:17:LEU:N	2.89	0.41
1:B:279:ARG:HD2	1:E:22:TRP:CG	2.55	0.41
1:D:86:LEU:HD11	1:F:35:LEU:O	2.21	0.41
1:D:401:LYS:NZ	4:D:602:HOH:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:36:GLN:HB3	1:E:44:SER:OG	2.21	0.41
1:F:223:SER:O	1:F:233:GLY:HA2	2.21	0.40
1:C:27:ARG:HD2	1:C:28:LEU:HG	2.04	0.40
1:F:144:ASP:OD1	1:F:145:LYS:N	2.55	0.40
1:F:305:HIS:CG	1:F:310:LEU:HD13	2.57	0.40
1:A:38:VAL:HG21	1:A:44:SER:HB3	2.03	0.40
1:A:151:MET:HE3	1:A:151:MET:HB2	1.96	0.40
1:B:272:ARG:O	1:B:276:ILE:HG13	2.22	0.40
1:B:356:TYR:OH	1:B:401:LYS:HE3	2.21	0.40
1:D:179:CYS:HA	1:D:185:VAL:HG21	2.01	0.40
3:D:501:PLR:H4A3	3:D:501:PLR:H5A1	1.88	0.40
1:E:292:THR:N	1:E:293:PRO:HD2	2.36	0.40
1:C:201:PRO:HB2	1:C:204:GLU:HB2	2.03	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	390/397 (98%)	373 (96%)	17 (4%)	0	100	100
1	B	393/397 (99%)	379 (96%)	13 (3%)	1 (0%)	36	66
1	C	391/397 (98%)	376 (96%)	15 (4%)	0	100	100
1	D	393/397 (99%)	385 (98%)	8 (2%)	0	100	100
1	E	389/397 (98%)	374 (96%)	15 (4%)	0	100	100
1	F	388/397 (98%)	376 (97%)	10 (3%)	2 (0%)	24	55
All	All	2344/2382 (98%)	2263 (96%)	78 (3%)	3 (0%)	48	77

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	92	THR
1	B	86	LEU
1	F	91	MET

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	313/316 (99%)	309 (99%)	4 (1%)	61	86
1	B	314/316 (99%)	301 (96%)	13 (4%)	27	62
1	C	312/316 (99%)	303 (97%)	9 (3%)	37	73
1	D	314/316 (99%)	300 (96%)	14 (4%)	24	58
1	E	310/316 (98%)	306 (99%)	4 (1%)	61	86
1	F	309/316 (98%)	300 (97%)	9 (3%)	37	73
All	All	1872/1896 (99%)	1819 (97%)	53 (3%)	38	73

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	14	LEU
1	C	17	LEU
1	C	86	LEU
1	C	90	ARG
1	C	95	LEU
1	C	127	VAL
1	C	157	LEU
1	C	226	VAL
1	C	310	LEU
1	A	91	MET
1	A	226	VAL
1	A	310	LEU
1	A	404	VAL
1	B	16	HIS
1	B	17	LEU
1	B	90	ARG

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Mol	Chain	Res	Type
1	B	127	VAL
1	B	140	VAL
1	B	157	LEU
1	B	208	LEU
1	B	220	GLU
1	B	223	SER
1	B	238	ARG
1	B	351	LEU
1	B	352	GLU
1	B	364	VAL
1	D	11	VAL
1	D	14	LEU
1	D	43	LEU
1	D	47	ASP
1	D	89	VAL
1	D	127	VAL
1	D	157	LEU
1	D	168	ARG
1	D	220	GLU
1	D	226	VAL
1	D	281	SER
1	D	310	LEU
1	D	364	VAL
1	D	402	LEU
1	E	31	HIS
1	E	127	VAL
1	E	190	ASP
1	E	198	THR
1	F	67	ARG
1	F	88	ARG
1	F	89	VAL
1	F	127	VAL
1	F	223	SER
1	F	244	ASP
1	F	281	SER
1	F	310	LEU
1	F	364	VAL

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

Mol	Chain	Res	Type
1	C	79	ASN

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Mol	Chain	Res	Type
1	C	115	ASN
1	C	335	ASN
1	B	79	ASN
1	B	115	ASN
1	D	31	HIS
1	D	79	ASN
1	E	36	GLN
1	E	333	GLN
1	E	335	ASN
1	F	36	GLN
1	F	360	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	PLR	E	501	-	15,15,15	1.13	2 (13%)	21,22,22	1.44	4 (19%)
2	PLP	B	501	1	15,15,16	1.17	2 (13%)	21,22,23	1.25	3 (14%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PLR	F	501	-	15,15,15	1.12	2 (13%)	21,22,22	1.33	3 (14%)
2	PLP	C	501	1	15,15,16	1.09	2 (13%)	21,22,23	1.04	2 (9%)
2	PLP	A	501	1	15,15,16	1.15	2 (13%)	21,22,23	1.19	1 (4%)
3	PLR	D	501	-	15,15,15	1.10	1 (6%)	21,22,22	1.43	4 (19%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PLR	E	501	-	-	3/6/6/6	0/1/1/1
2	PLP	B	501	1	-	3/6/6/8	0/1/1/1
3	PLR	F	501	-	-	0/6/6/6	0/1/1/1
2	PLP	C	501	1	-	3/6/6/8	0/1/1/1
2	PLP	A	501	1	-	3/6/6/8	0/1/1/1
3	PLR	D	501	-	-	3/6/6/6	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	PLP	C2-N1	2.73	1.38	1.33
3	E	501	PLR	C2-N1	2.66	1.38	1.33
2	A	501	PLP	C2-N1	2.62	1.38	1.33
2	C	501	PLP	C2-N1	2.62	1.38	1.33
3	F	501	PLR	C2-N1	2.60	1.38	1.33
3	D	501	PLR	C2-N1	2.59	1.38	1.33
2	A	501	PLP	C6-N1	2.18	1.38	1.34
2	B	501	PLP	C6-N1	2.12	1.38	1.34
3	F	501	PLR	C6-N1	2.08	1.38	1.34
3	E	501	PLR	C6-N1	2.07	1.38	1.34
2	C	501	PLP	C6-N1	2.05	1.38	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	501	PLR	C4A-C4-C5	-3.46	117.37	120.94
3	D	501	PLR	C6-C5-C4	3.17	120.69	118.10
3	E	501	PLR	C4A-C4-C5	-3.08	117.77	120.94
3	E	501	PLR	C6-C5-C4	2.86	120.44	118.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	501	PLR	C4A-C4-C5	-2.80	118.05	120.94
3	F	501	PLR	C6-C5-C4	2.54	120.18	118.10
2	B	501	PLP	C5-C6-N1	-2.48	119.79	123.83
3	E	501	PLR	C5-C6-N1	-2.48	119.81	123.83
2	B	501	PLP	C6-C5-C4	2.46	120.12	118.10
3	D	501	PLR	C5-C6-N1	-2.43	119.87	123.83
3	E	501	PLR	C5A-C5-C4	-2.32	118.05	122.64
3	F	501	PLR	C5-C6-N1	-2.32	120.06	123.83
2	B	501	PLP	C5A-C5-C4	-2.26	118.17	122.64
2	C	501	PLP	C5-C6-N1	-2.26	120.16	123.83
2	A	501	PLP	C5-C6-N1	-2.18	120.29	123.83
2	C	501	PLP	C6-C5-C4	2.11	119.83	118.10
3	D	501	PLR	C5A-C5-C4	-2.05	118.59	122.64

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	501	PLP	C5A-O4P-P-O2P
2	C	501	PLP	C5A-O4P-P-O3P
2	A	501	PLP	C5A-O4P-P-O1P
2	A	501	PLP	C5A-O4P-P-O2P
2	A	501	PLP	C5A-O4P-P-O3P
2	B	501	PLP	C5A-O4P-P-O1P
2	B	501	PLP	C5A-O4P-P-O2P
2	B	501	PLP	C5A-O4P-P-O3P
3	D	501	PLR	C5A-O4P-P-O2P
3	E	501	PLR	C5A-O4P-P-O1P
3	E	501	PLR	C5A-O4P-P-O2P
3	E	501	PLR	C5A-O4P-P-O3P
3	D	501	PLR	C5A-O4P-P-O1P
3	D	501	PLR	C5A-O4P-P-O3P
2	C	501	PLP	C5A-O4P-P-O1P

There are no ring outliers.

4 monomers are involved in 9 short contacts:

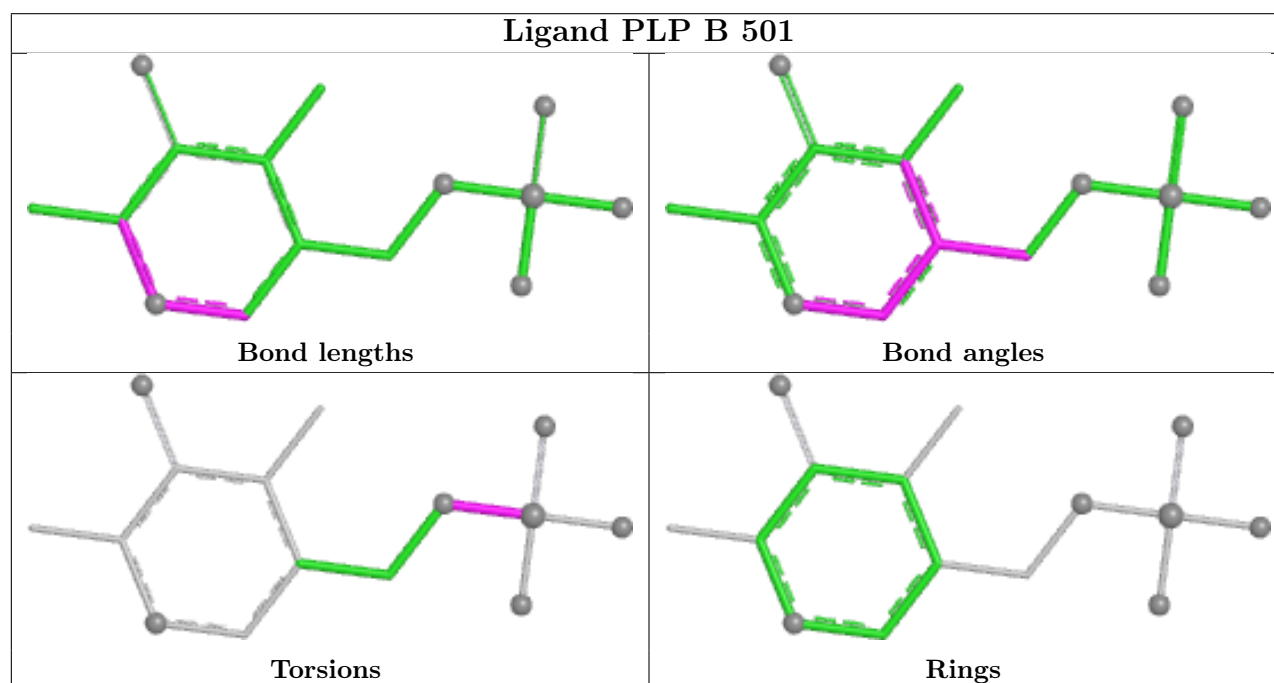
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	501	PLR	2	0
2	B	501	PLP	2	0
3	F	501	PLR	4	0

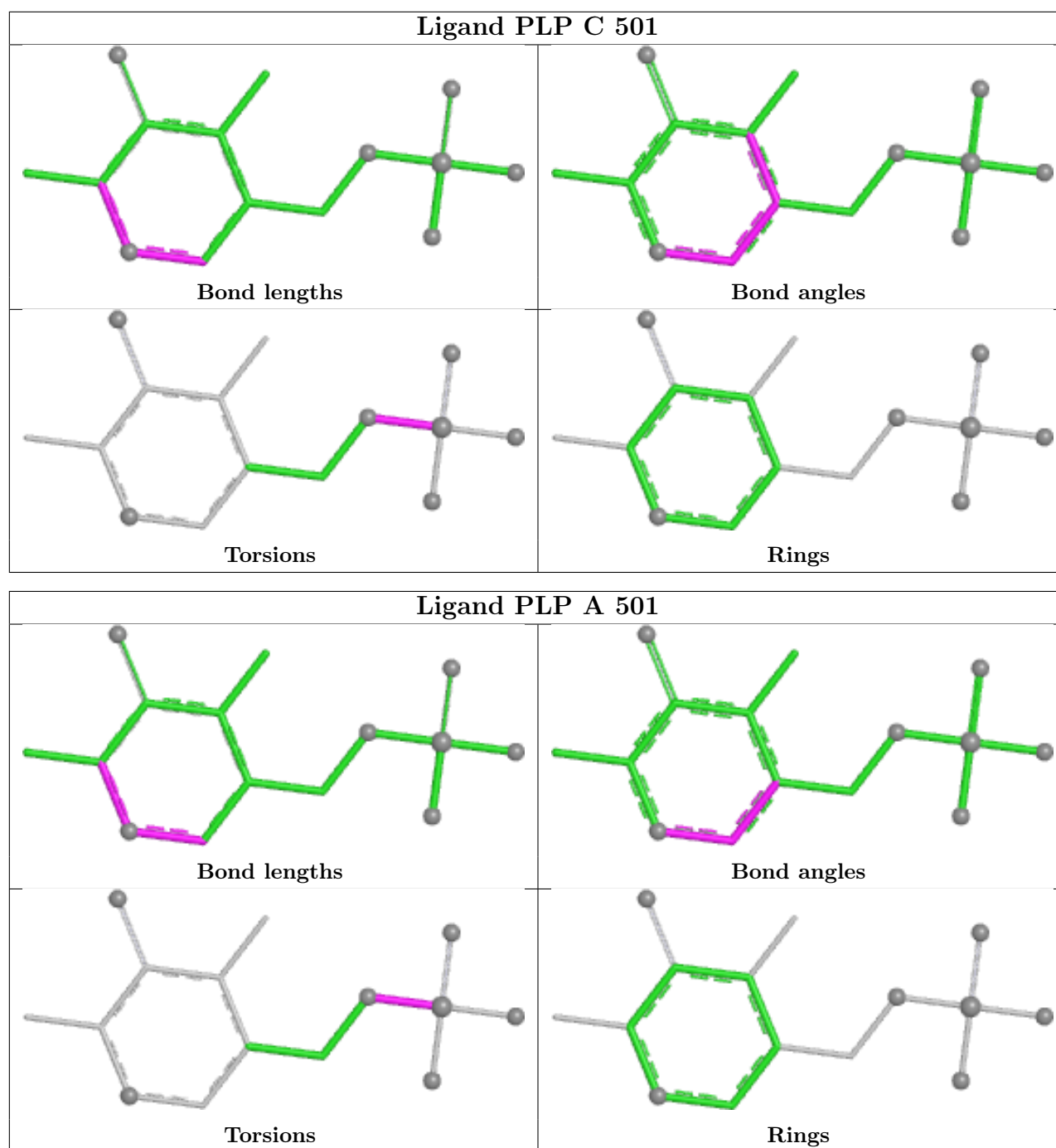
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	501	PLR	1	0

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	394/397 (99%)	-0.09	5 (1%) 75 66	37, 62, 96, 140	0
1	B	395/397 (99%)	-0.14	4 (1%) 79 72	38, 65, 96, 136	0
1	C	393/397 (98%)	-0.01	9 (2%) 61 51	38, 66, 124, 184	0
1	D	395/397 (99%)	-0.22	5 (1%) 75 66	42, 60, 87, 130	0
1	E	391/397 (98%)	0.02	4 (1%) 79 72	42, 70, 105, 134	0
1	F	390/397 (98%)	-0.12	6 (1%) 72 63	44, 66, 93, 144	0
All	All	2358/2382 (98%)	-0.09	33 (1%) 73 64	37, 65, 100, 184	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	87	GLY	5.0
1	F	91	MET	4.9
1	A	87	GLY	4.0
1	C	14	LEU	4.0
1	F	89	VAL	3.7
1	C	30	VAL	3.5
1	D	86	LEU	3.4
1	F	86	LEU	3.4
1	A	88	ARG	3.4
1	A	86	LEU	3.2
1	E	86	LEU	3.1
1	C	87	GLY	3.1
1	B	14	LEU	3.1
1	A	14	LEU	3.0
1	E	87	GLY	2.9
1	C	86	LEU	2.8
1	B	86	LEU	2.8
1	C	24	SER	2.7
1	E	14	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	F	17	LEU	2.6
1	D	87	GLY	2.5
1	D	11	VAL	2.4
1	F	87	GLY	2.3
1	B	84	SER	2.3
1	D	404	VAL	2.3
1	C	18	SER	2.3
1	C	29	GLY	2.3
1	A	90	ARG	2.2
1	E	224	THR	2.2
1	F	197	GLY	2.2
1	C	88	ARG	2.1
1	C	17	LEU	2.1
1	D	286	TRP	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.

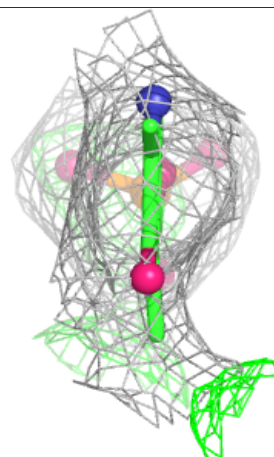
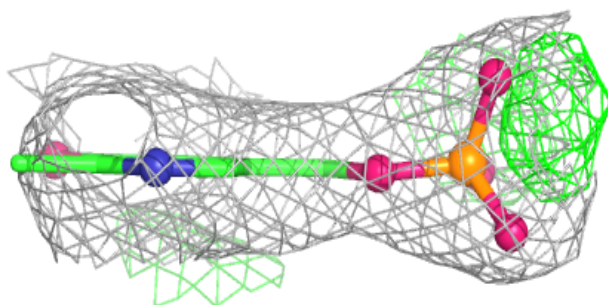
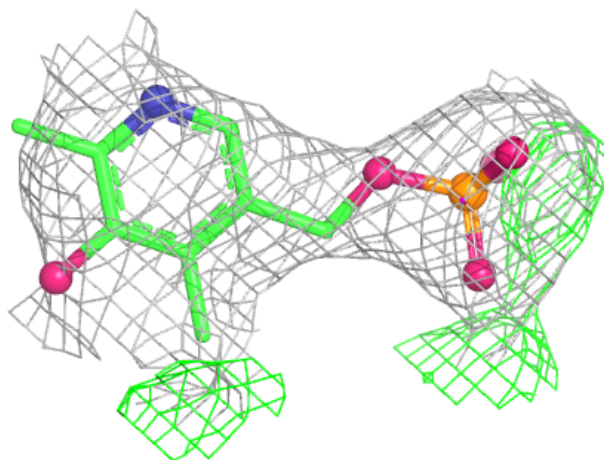
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PLP	A	501	15/16	0.86	0.12	48,69,88,89	0
2	PLP	C	501	15/16	0.92	0.09	50,67,73,76	0
3	PLR	F	501	15/15	0.92	0.10	51,64,78,84	0
3	PLR	D	501	15/15	0.93	0.09	50,57,65,75	0
3	PLR	E	501	15/15	0.93	0.09	58,70,84,85	0
2	PLP	B	501	15/16	0.93	0.09	51,59,75,81	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.

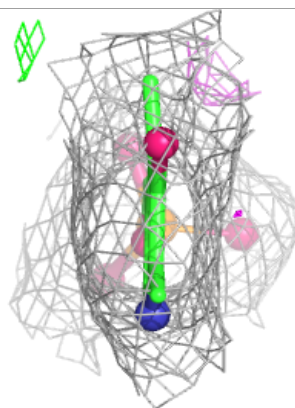
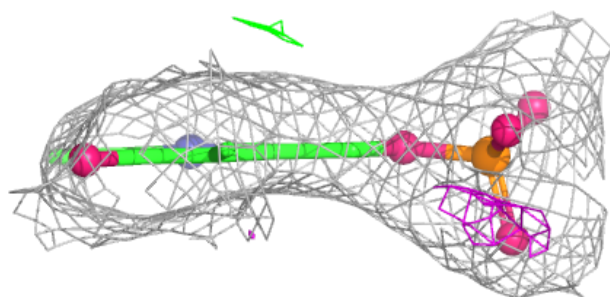
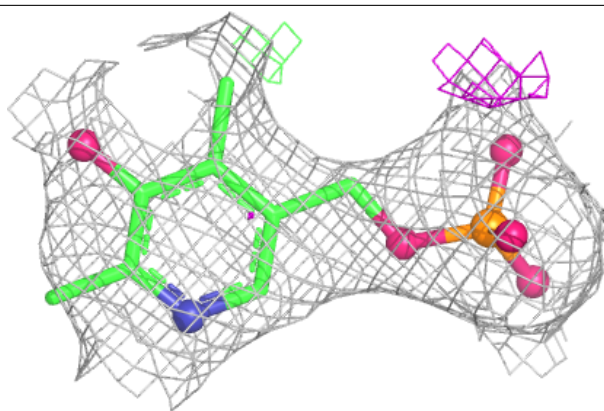
Electron density around PLP A 501:

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

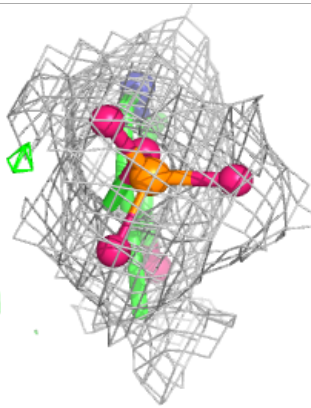
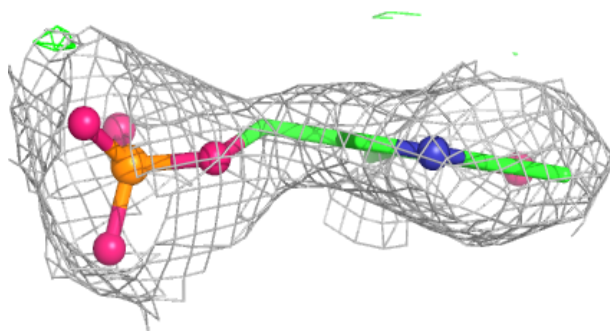
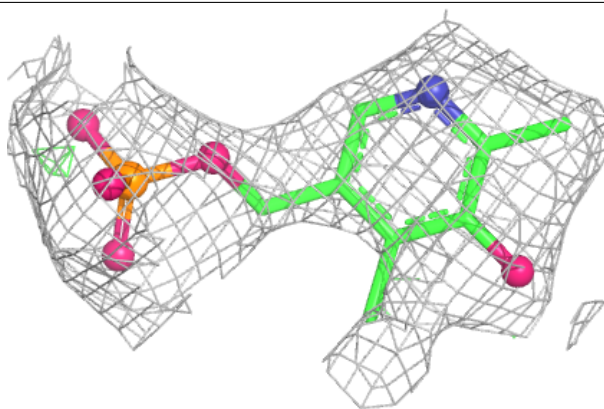


Electron density around PLP C 501:

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

**Electron density around PLP B 501:**

$2mF_o-DF_c$ (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.