



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:34 AM UTC

PDB ID : 3RBF / pdb_00003rbf
Title : Crystal structure of Human aromatic L-amino acid decarboxylase (AADC) in the apo form
Authors : Giardina, G.; Montioli, R.; Gianni, S.; Cellini, B.; Paiardini, A.; Borri Voltattorni, C.; Cutruzzola, F.
Deposited on : 2011-03-29
Resolution : 2.90 Å(reported)

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

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

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
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

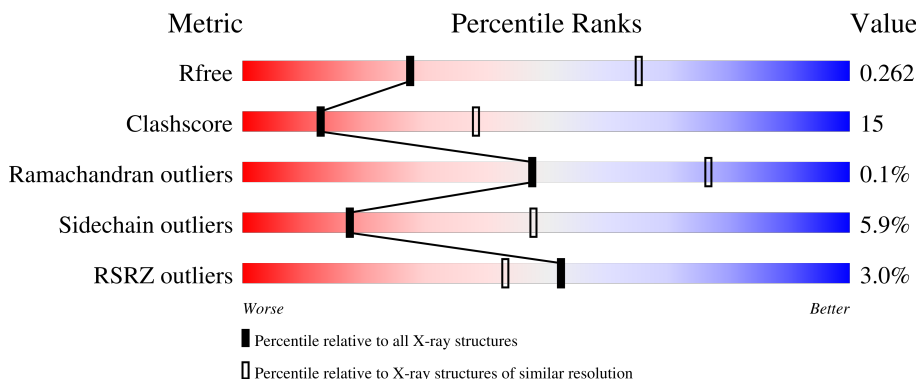
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	480	
1	B	480	

2 Entry composition [i](#)

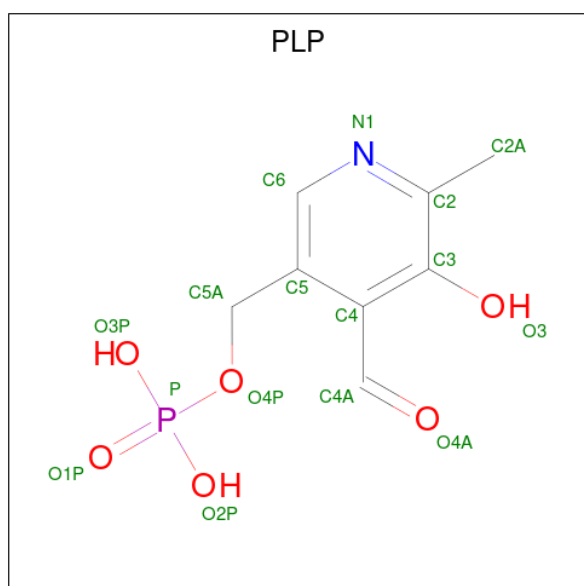
There are 3 unique types of molecules in this entry. The entry contains 6972 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Aromatic-L-amino-acid decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	443	Total	C	N	O	S	0	0	0
			3474	2224	600	623	27			
1	B	443	Total	C	N	O	S	0	0	0
			3480	2229	601	624	26			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

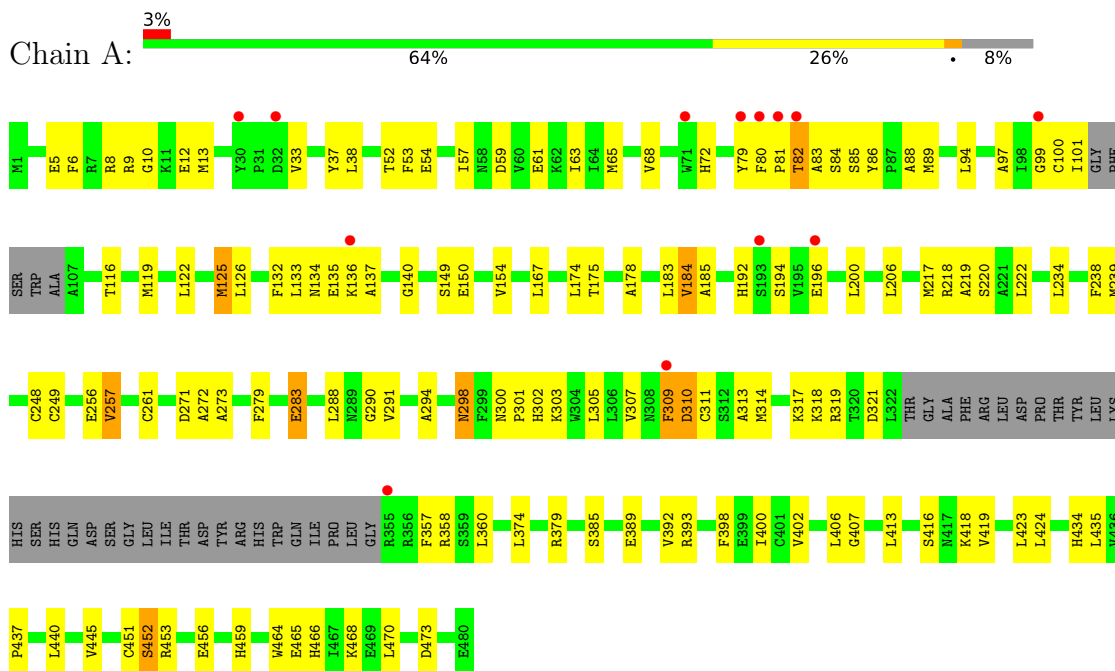
- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Cl	0	0
			2	2		

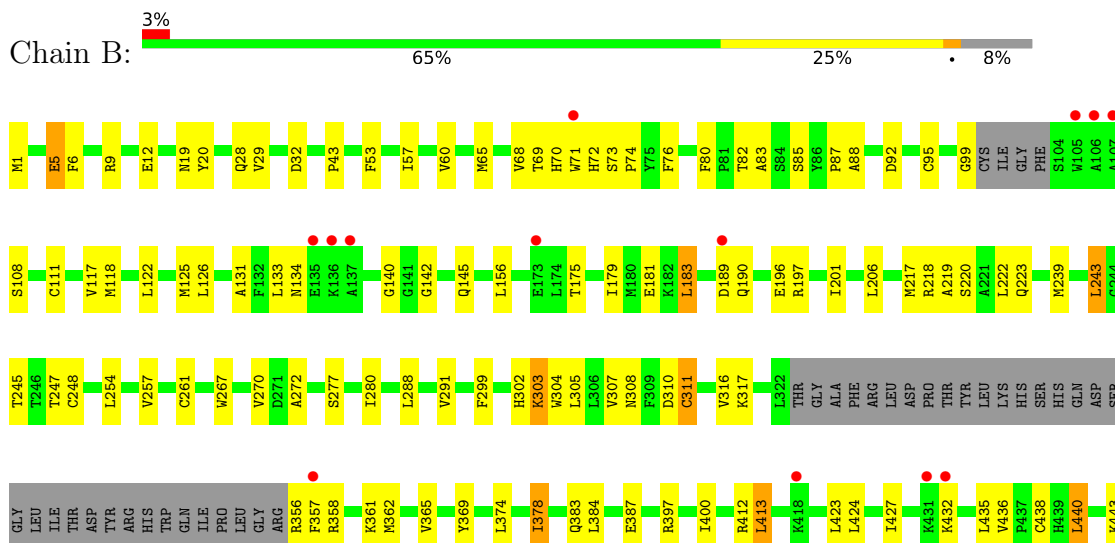
3 Residue-property plots

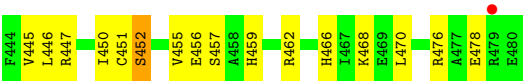
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Aromatic-L-amino-acid decarboxylase



• Molecule 1: Aromatic-L-amino-acid decarboxylase





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	175.84Å 175.84Å 74.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.77 – 2.90 28.77 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.1 (28.77-2.90) 98.9 (28.77-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.10 (at 2.91Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.212 , 0.267 0.209 , 0.262	Depositor DCC
R_{free} test set	1338 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	45.5	Xtriage
Anisotropy	0.405	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 33.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6972	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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: CL, 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	0.67	0/3553	0.97	6/4805 (0.1%)
1	B	0.69	0/3561	0.95	3/4817 (0.1%)
All	All	0.68	0/7114	0.96	9/9622 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	86	TYR	CA-C-N	-6.90	112.67	119.64
1	A	86	TYR	C-N-CA	-6.90	112.67	119.64
1	A	65	MET	CA-C-N	-6.52	112.97	119.56
1	A	65	MET	C-N-CA	-6.52	112.97	119.56
1	A	309	PHE	N-CA-C	5.97	118.55	111.33
1	B	311	CYS	N-CA-C	5.79	117.97	108.52
1	B	378	ILE	CB-CA-C	-5.24	105.17	111.88
1	A	434	HIS	N-CA-C	5.08	117.68	108.69
1	B	303	LYS	N-CA-C	5.02	116.83	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	3474	0	3464	106	0
1	B	3480	0	3466	109	0
2	A	16	0	8	5	0
3	B	2	0	0	0	0
All	All	6972	0	6938	209	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 15.

All (209) 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:53:PHE:CE2	1:A:57:ILE:HD11	1.81	1.14
1:A:302:HIS:ND1	1:A:309:PHE:O	1.87	1.06
1:A:83:ALA:HB3	1:A:302:HIS:O	1.58	1.03
1:B:412:ARG:HD3	1:B:443:LYS:HE2	1.39	1.02
1:B:304:TRP:HB3	1:B:378:ILE:CD1	1.93	0.98
1:B:189:ASP:OD1	1:B:190:GLN:HG2	1.67	0.93
1:A:217:MET:SD	1:A:239:MET:HE1	2.12	0.89
1:A:456:GLU:H	1:A:459:HIS:HD2	1.19	0.89
1:B:239:MET:CE	1:B:261:CYS:SG	2.64	0.85
1:B:456:GLU:H	1:B:459:HIS:HD2	1.25	0.85
1:B:53:PHE:CE2	1:B:57:ILE:HD11	2.12	0.85
1:B:239:MET:HE2	1:B:261:CYS:SG	2.16	0.84
1:B:304:TRP:HB3	1:B:378:ILE:HD11	1.57	0.83
1:B:288:LEU:CD1	1:B:291:VAL:HG21	2.10	0.81
1:A:38:LEU:HD11	1:A:63:ILE:HG22	1.62	0.80
1:B:1:MET:HE3	1:B:5:GLU:HB3	1.63	0.80
1:A:288:LEU:O	1:A:291:VAL:HG23	1.81	0.79
1:B:288:LEU:HD11	1:B:291:VAL:HG21	1.65	0.78
1:A:101:ILE:HB	1:B:82:THR:HG21	1.66	0.75
1:B:95:CYS:SG	1:B:357:PHE:HB2	2.27	0.74
1:B:412:ARG:CD	1:B:443:LYS:HE2	2.17	0.74
1:B:369:TYR:CB	1:B:374:LEU:CD1	2.66	0.74
1:A:288:LEU:HG	1:A:291:VAL:CG2	2.16	0.73
1:B:307:VAL:HG11	1:B:362:MET:HG2	1.70	0.73
1:A:38:LEU:HD11	1:A:63:ILE:CG2	2.18	0.73
1:A:456:GLU:H	1:A:459:HIS:CD2	2.05	0.73
1:B:432:LYS:HE2	1:B:462:ARG:NH1	2.04	0.73
1:B:412:ARG:HD3	1:B:443:LYS:CE	2.19	0.72
1:A:100:CYS:O	1:A:101:ILE:HG13	1.89	0.72
1:A:83:ALA:CB	1:A:302:HIS:O	2.36	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:ARG:O	1:B:219:ALA:HB3	1.89	0.70
1:B:305:LEU:HD22	1:B:374:LEU:HD23	1.73	0.70
1:A:38:LEU:CD1	1:A:63:ILE:CG2	2.70	0.70
1:A:303:LYS:NZ	2:A:481:PLP:C4A	2.56	0.69
1:B:189:ASP:OD1	1:B:190:GLN:N	2.26	0.68
1:B:369:TYR:HB2	1:B:374:LEU:CD1	2.23	0.68
1:B:29:VAL:O	1:B:69:THR:O	2.12	0.68
1:B:28:GLN:NE2	1:B:32:ASP:HB2	2.09	0.67
1:A:389:GLU:OE2	1:A:393:ARG:NH2	2.27	0.67
1:B:196:GLU:HA	1:B:206:LEU:HD22	1.76	0.66
1:A:398:PHE:CE2	1:A:413:LEU:HD13	2.31	0.66
1:B:272:ALA:HB3	1:B:299:PHE:HB3	1.76	0.66
1:A:79:TYR:HA	1:A:451:CYS:SG	2.36	0.65
1:A:302:HIS:CE1	1:A:309:PHE:O	2.49	0.64
1:A:416:SER:OG	1:A:419:VAL:HG23	1.97	0.64
1:B:134:ASN:OD1	1:B:142:GLY:N	2.30	0.63
1:A:5:GLU:OE2	1:A:9:ARG:NE	2.31	0.63
1:A:59:ASP:HB3	1:A:63:ILE:HD12	1.81	0.63
1:A:99:GLY:HA3	1:B:80:PHE:HZ	1.64	0.63
1:B:71:TRP:HE3	1:B:76:PHE:CZ	2.17	0.63
1:B:71:TRP:CE3	1:B:76:PHE:CZ	2.87	0.63
1:B:307:VAL:HG12	1:B:308:ASN:H	1.62	0.62
1:A:38:LEU:CD1	1:A:63:ILE:HG22	2.29	0.62
1:A:192:HIS:CD2	1:A:194:SER:H	2.17	0.62
1:A:8:ARG:NH2	1:A:9:ARG:HE	1.97	0.62
1:A:192:HIS:HD2	1:A:194:SER:H	1.47	0.62
1:B:302:HIS:HD2	1:B:307:VAL:O	1.81	0.62
1:B:356:ARG:O	1:B:358:ARG:HG3	1.98	0.62
1:B:387:GLU:OE2	1:B:457:SER:HB2	2.00	0.61
1:A:288:LEU:HG	1:A:291:VAL:HG21	1.83	0.60
1:A:456:GLU:N	1:A:459:HIS:HD2	1.94	0.60
1:A:303:LYS:NZ	2:A:481:PLP:H4A	2.15	0.59
1:B:196:GLU:HA	1:B:206:LEU:CD2	2.31	0.59
1:A:5:GLU:OE2	1:A:9:ARG:CZ	2.50	0.59
1:B:53:PHE:CZ	1:B:57:ILE:HD11	2.38	0.59
1:A:303:LYS:HZ2	2:A:481:PLP:C4A	2.16	0.58
1:A:38:LEU:HD12	1:A:63:ILE:HG23	1.85	0.58
1:B:219:ALA:HA	1:B:222:LEU:HB3	1.85	0.57
1:B:456:GLU:H	1:B:459:HIS:CD2	2.13	0.57
1:A:119:MET:HE1	1:A:313:ALA:HB1	1.85	0.57
1:B:476:ARG:HG3	1:B:476:ARG:HH11	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLU:HA	1:A:206:LEU:HD22	1.86	0.57
1:B:369:TYR:HB3	1:B:374:LEU:CD1	2.35	0.57
1:B:452:SER:OG	1:B:455:VAL:HG23	2.03	0.57
1:B:369:TYR:CB	1:B:374:LEU:HD12	2.35	0.56
1:A:134:ASN:HB3	1:A:140:GLY:O	2.04	0.56
1:A:175:THR:HG23	1:A:178:ALA:H	1.70	0.56
1:B:85:SER:HB2	1:B:87:PRO:HD2	1.87	0.56
1:A:122:LEU:HA	1:A:125:MET:HE3	1.87	0.56
1:A:239:MET:HG2	1:A:261:CYS:SG	2.46	0.56
1:A:301:PRO:HD2	1:A:311:CYS:O	2.06	0.56
1:B:69:THR:O	1:B:70:HIS:HB2	2.05	0.56
1:A:294:ALA:O	1:A:317:LYS:HE2	2.05	0.56
1:B:369:TYR:HB3	1:B:374:LEU:HD12	1.87	0.56
1:A:150:GLU:O	1:A:154:VAL:HG23	2.05	0.56
1:A:99:GLY:HA3	1:B:80:PHE:CZ	2.42	0.55
1:B:71:TRP:HB2	1:B:80:PHE:CZ	2.42	0.55
1:A:100:CYS:O	1:A:101:ILE:CG1	2.56	0.54
1:B:456:GLU:N	1:B:459:HIS:HD2	2.02	0.54
1:A:126:LEU:HD22	1:A:279:PHE:HB2	1.91	0.53
1:A:300:ASN:ND2	2:A:481:PLP:O3P	2.41	0.53
1:A:357:PHE:CE1	1:A:360:LEU:HB2	2.44	0.53
1:A:470:LEU:O	1:A:473:ASP:HB2	2.09	0.53
1:A:184:VAL:CG2	1:A:234:LEU:HB3	2.39	0.52
1:B:423:LEU:HD23	1:B:446:LEU:HD11	1.91	0.52
1:B:438:CYS:SG	1:B:447:ARG:HD2	2.50	0.52
1:A:318:LYS:O	1:A:321:ASP:HB2	2.10	0.52
1:B:308:ASN:HB3	1:B:361:LYS:HD2	1.91	0.52
1:A:94:LEU:HD23	1:A:360:LEU:HD21	1.92	0.52
1:A:440:LEU:HD12	1:A:445:VAL:HG21	1.92	0.52
1:B:277:SER:HB2	1:B:304:TRP:HB2	1.92	0.52
1:A:5:GLU:OE2	1:A:9:ARG:NH2	2.43	0.51
1:A:302:HIS:CG	1:A:309:PHE:O	2.63	0.51
1:A:310:ASP:CG	1:A:310:ASP:O	2.54	0.51
1:A:94:LEU:HD21	1:B:60:VAL:HG11	1.92	0.51
1:B:308:ASN:OD1	1:B:310:ASP:HB3	2.10	0.51
1:A:136:LYS:O	1:A:137:ALA:HB3	2.11	0.51
1:A:218:ARG:NH1	1:A:256:GLU:OE2	2.43	0.51
1:B:239:MET:HE1	1:B:257:VAL:HG13	1.91	0.51
1:A:302:HIS:HB2	1:A:309:PHE:HA	1.93	0.50
1:B:369:TYR:HB2	1:B:374:LEU:HD13	1.93	0.50
1:A:38:LEU:HD12	1:A:63:ILE:CG2	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:ALA:O	1:B:223:GLN:N	2.34	0.50
1:B:108:SER:HB3	1:B:111:CYS:SG	2.52	0.50
1:B:440:LEU:HD22	1:B:445:VAL:HG21	1.92	0.50
1:B:189:ASP:CG	1:B:190:GLN:HE21	2.20	0.49
1:B:427:ILE:HG21	1:B:435:LEU:HD13	1.94	0.49
1:B:239:MET:HE1	1:B:261:CYS:SG	2.50	0.49
1:B:310:ASP:OD1	1:B:310:ASP:O	2.30	0.49
1:A:219:ALA:O	1:A:220:SER:C	2.55	0.49
1:A:357:PHE:CD1	1:A:360:LEU:HB2	2.48	0.49
1:A:303:LYS:HZ1	2:A:481:PLP:H4A	1.77	0.48
1:A:248:CYS:HB2	1:A:402:VAL:HG21	1.96	0.48
1:B:267:TRP:CD1	1:B:267:TRP:C	2.91	0.48
1:A:6:PHE:O	1:A:10:GLY:N	2.41	0.48
1:A:97:ALA:O	1:B:72:HIS:HD2	1.96	0.48
1:B:243:LEU:HD22	1:B:270:VAL:HG11	1.95	0.48
1:B:303:LYS:HE3	1:B:451:CYS:SG	2.54	0.48
1:A:192:HIS:HD2	1:A:194:SER:HB3	1.78	0.47
1:A:53:PHE:CD2	1:A:57:ILE:HD11	2.43	0.47
1:A:33:VAL:HG11	1:A:37:TYR:CG	2.49	0.47
1:B:1:MET:HE2	1:B:6:PHE:HA	1.95	0.47
1:B:218:ARG:HG3	1:B:218:ARG:HH11	1.80	0.47
1:B:304:TRP:HB3	1:B:378:ILE:HD12	1.89	0.47
1:A:33:VAL:HG11	1:A:37:TYR:CD2	2.50	0.47
1:B:126:LEU:HD11	1:B:299:PHE:CE2	2.49	0.47
1:B:243:LEU:HD22	1:B:270:VAL:CG1	2.45	0.47
1:A:406:LEU:O	1:A:406:LEU:HG	2.14	0.46
1:A:305:LEU:HD22	1:A:374:LEU:HD22	1.97	0.46
1:B:218:ARG:HG3	1:B:218:ARG:NH1	2.31	0.46
1:B:288:LEU:CD1	1:B:291:VAL:CG2	2.90	0.46
1:A:72:HIS:NE2	1:A:84:SER:HB2	2.31	0.46
1:B:369:TYR:CB	1:B:374:LEU:HD13	2.45	0.46
1:A:79:TYR:CZ	1:A:81:PRO:HA	2.51	0.45
1:B:88:ALA:O	1:B:92:ASP:HB2	2.16	0.45
1:B:452:SER:OG	1:B:455:VAL:CG2	2.65	0.45
1:A:12:GLU:OE1	1:A:53:PHE:HD2	1.99	0.45
1:A:283:GLU:OE2	1:A:379:ARG:NH2	2.43	0.45
1:A:413:LEU:HD22	1:A:423:LEU:HD22	1.99	0.45
1:A:271:ASP:OD1	1:A:271:ASP:C	2.60	0.45
1:A:300:ASN:HB2	1:A:303:LYS:HG3	1.99	0.45
1:B:310:ASP:OD2	1:B:358:ARG:HD3	2.17	0.45
1:B:197:ARG:HE	1:B:201:ILE:HD11	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:ALA:O	1:B:72:HIS:CD2	2.69	0.45
1:B:99:GLY:O	1:B:357:PHE:CE2	2.70	0.45
1:A:82:THR:HA	1:A:309:PHE:HB3	1.99	0.44
1:A:272:ALA:O	1:A:273:ALA:C	2.59	0.44
1:B:436:VAL:O	1:B:447:ARG:HB2	2.17	0.44
1:A:298:ASN:HB2	1:A:314:MET:HG3	2.00	0.44
1:B:19:ASN:O	1:B:20:TYR:C	2.60	0.44
1:B:74:PRO:HB3	1:B:455:VAL:HA	1.99	0.44
1:B:117:VAL:HG12	1:B:118:MET:HE2	1.98	0.44
1:A:465:GLU:O	1:A:468:LYS:HG2	2.17	0.44
1:B:80:PHE:CD2	1:B:83:ALA:HB2	2.53	0.44
1:B:122:LEU:HA	1:B:125:MET:HE3	1.99	0.44
1:A:385:SER:HB2	1:A:407:GLY:HA2	1.99	0.44
1:A:132:PHE:C	1:A:133:LEU:HD12	2.43	0.44
1:A:290:GLY:O	1:A:291:VAL:C	2.61	0.44
1:B:308:ASN:H	1:B:308:ASN:HD22	1.66	0.44
1:B:87:PRO:HB2	1:B:365:VAL:HG22	2.00	0.44
1:B:307:VAL:HG12	1:B:308:ASN:N	2.31	0.43
1:A:13:MET:HE2	1:A:53:PHE:HZ	1.83	0.43
1:A:288:LEU:HG	1:A:291:VAL:HG23	2.00	0.43
1:A:52:THR:HG22	1:A:54:GLU:OE1	2.18	0.43
1:A:134:ASN:HD22	1:A:319:ARG:HH21	1.65	0.43
1:B:245:THR:OG1	1:B:248:CYS:O	2.33	0.43
1:B:277:SER:CB	1:B:304:TRP:HB2	2.48	0.43
1:A:452:SER:O	1:A:453:ARG:C	2.61	0.43
1:B:28:GLN:HE21	1:B:32:ASP:HB2	1.80	0.43
1:A:85:SER:O	1:A:88:ALA:HB3	2.19	0.43
1:A:392:VAL:HG22	1:A:464:TRP:CZ3	2.54	0.43
1:B:85:SER:CB	1:B:87:PRO:HD2	2.50	0.42
1:A:248:CYS:HB3	1:A:440:LEU:HD21	2.02	0.42
1:A:249:CYS:SG	1:A:406:LEU:CD2	3.08	0.42
1:A:466:HIS:O	1:A:470:LEU:HG	2.20	0.42
1:B:397:ARG:O	1:B:413:LEU:HD12	2.20	0.42
1:B:9:ARG:HD2	1:B:12:GLU:OE1	2.20	0.41
1:A:85:SER:O	1:A:89:MET:HG3	2.20	0.41
1:A:302:HIS:HA	1:A:307:VAL:O	2.20	0.41
1:B:310:ASP:CG	1:B:358:ARG:HD3	2.45	0.41
1:B:384:LEU:CD2	1:B:457:SER:HB3	2.50	0.41
1:A:149:SER:HA	1:A:194:SER:HB2	2.03	0.41
1:A:174:LEU:HD12	1:A:174:LEU:HA	1.95	0.41
1:A:222:LEU:HD22	1:A:257:VAL:HG22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:CYS:SG	1:A:406:LEU:HD21	2.60	0.41
1:A:424:LEU:HD22	1:A:437:PRO:HG3	2.03	0.41
1:B:69:THR:O	1:B:70:HIS:CB	2.67	0.41
1:B:179:ILE:O	1:B:183:LEU:HD13	2.20	0.41
1:B:357:PHE:N	1:B:357:PHE:CD1	2.88	0.41
1:B:450:ILE:HG23	1:B:455:VAL:HG11	2.02	0.41
1:B:125:MET:HB3	1:B:280:ILE:HG22	2.03	0.40
1:B:254:LEU:HD12	1:B:254:LEU:HA	1.92	0.40
1:A:185:ALA:CB	1:A:238:PHE:HB3	2.51	0.40
1:B:218:ARG:HD2	1:B:218:ARG:HA	1.78	0.40
1:B:131:ALA:O	1:B:140:GLY:HA3	2.22	0.40
1:A:300:ASN:HA	1:A:301:PRO:HD3	1.96	0.40
1:B:60:VAL:O	1:B:65:MET:HG2	2.22	0.40
1:B:156:LEU:HD23	1:B:156:LEU:C	2.46	0.40
1:B:466:HIS:CE1	1:B:470:LEU:HD11	2.55	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	437/480 (91%)	411 (94%)	26 (6%)	0	100	100
1	B	437/480 (91%)	410 (94%)	26 (6%)	1 (0%)	43	72
All	All	874/960 (91%)	821 (94%)	52 (6%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	43	PRO

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	366/399 (92%)	346 (94%)	20 (6%)	19	50
1	B	366/399 (92%)	343 (94%)	23 (6%)	16	45
All	All	732/798 (92%)	689 (94%)	43 (6%)	18	48

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	GLU
1	A	68	VAL
1	A	80	PHE
1	A	82	THR
1	A	116	THR
1	A	125	MET
1	A	135	GLU
1	A	167	LEU
1	A	183	LEU
1	A	184	VAL
1	A	200	LEU
1	A	257	VAL
1	A	283	GLU
1	A	298	ASN
1	A	310	ASP
1	A	358	ARG
1	A	400	ILE
1	A	418	LYS
1	A	435	LEU
1	A	452	SER
1	B	5	GLU
1	B	68	VAL
1	B	73	SER
1	B	133	LEU
1	B	145	GLN
1	B	175	THR
1	B	181	GLU

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Mol	Chain	Res	Type
1	B	183	LEU
1	B	217	MET
1	B	220	SER
1	B	243	LEU
1	B	247	THR
1	B	311	CYS
1	B	316	VAL
1	B	317	LYS
1	B	383	GLN
1	B	400	ILE
1	B	413	LEU
1	B	424	LEU
1	B	440	LEU
1	B	452	SER
1	B	468	LYS
1	B	478	GLU

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	19	ASN
1	A	48	GLN
1	A	176	GLN
1	A	192	HIS
1	A	298	ASN
1	A	375	GLN
1	A	381	HIS
1	A	383	GLN
1	A	434	HIS
1	A	459	HIS
1	B	2	ASN
1	B	28	GLN
1	B	72	HIS
1	B	145	GLN
1	B	298	ASN
1	B	308	ASN
1	B	386	HIS
1	B	434	HIS
1	B	459	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PLP	A	481	-	16,16,16	1.31	2 (12%)	20,23,23	1.34	4 (20%)

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	PLP	A	481	-	-	4/8/8/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	481	PLP	C2-N1	2.87	1.39	1.33
2	A	481	PLP	C6-N1	2.16	1.38	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	481	PLP	O4A-C4A-C4	-2.91	117.90	124.80
2	A	481	PLP	C3-C4-C5	2.60	120.36	118.28
2	A	481	PLP	O3P-P-O4P	2.55	113.32	106.67
2	A	481	PLP	C5-C6-N1	-2.13	120.37	123.83

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	481	PLP	C5A-O4P-P-O1P
2	A	481	PLP	C5A-O4P-P-O3P
2	A	481	PLP	C3-C4-C4A-O4A
2	A	481	PLP	C5A-O4P-P-O2P

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	481	PLP	5	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	443/480 (92%)	-0.14	13 (2%)	53	45	21, 37, 63, 80	5 (1%)
1	B	443/480 (92%)	-0.11	14 (3%)	50	41	20, 39, 61, 86	6 (1%)
All	All	886/960 (92%)	-0.12	27 (3%)	52	43	20, 38, 62, 86	11 (1%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	80	PHE	8.7
1	A	309	PHE	7.4
1	B	136	LYS	7.1
1	B	431	LYS	6.4
1	B	357	PHE	6.4
1	B	418	LYS	4.8
1	A	81	PRO	4.3
1	B	107	ALA	4.0
1	B	479	ARG	3.5
1	B	137	ALA	3.4
1	B	189	ASP	3.2
1	B	106	ALA	3.0
1	A	196	GLU	3.0
1	B	105	TRP	3.0
1	A	79	TYR	2.8
1	B	135	GLU	2.6
1	A	82	THR	2.5
1	A	71	TRP	2.5
1	A	99	GLY	2.5
1	B	71	TRP	2.3
1	A	136	LYS	2.3
1	A	32	ASP	2.3
1	B	432	LYS	2.2
1	A	355	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	193	SER	2.1
1	B	173	GLU	2.1
1	A	30	TYR	2.1

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

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

6.3 Carbohydrates [i](#)

There are no 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
2	PLP	A	481	16/16	0.77	0.25	37,39,41,44	16
3	CL	B	482	1/1	0.92	0.08	45,45,45,45	0
3	CL	B	481	1/1	0.97	0.04	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.