



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 8, 2025 – 02:09 PM JST

PDB ID : 9JYX / pdb_00009jyx
Title : Crystal structure of Nir2 DDHD domain
Authors : Kim, D.; Lee, C.
Deposited on : 2024-10-13
Resolution : 2.80 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
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) : 2.46

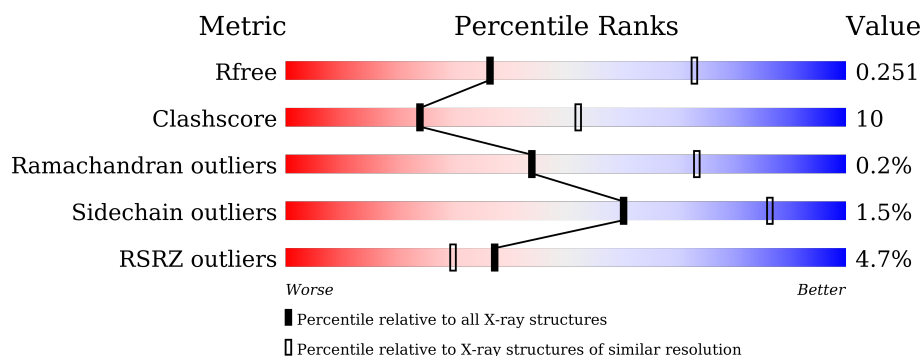
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	164625	3657 (2.80-2.80)
Clashscore	180529	4123 (2.80-2.80)
Ramachandran outliers	177936	4071 (2.80-2.80)
Sidechain outliers	177891	4073 (2.80-2.80)
RSRZ outliers	164620	3659 (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	360	<div> <div>3%</div> <div> <div></div> <div>70%</div> <div>14%</div> <div>•</div> <div>14%</div> </div> </div>
1	B	360	<div> <div>3%</div> <div> <div></div> <div>66%</div> <div>18%</div> <div>•</div> <div>16%</div> </div> </div>
1	C	360	<div> <div>6%</div> <div> <div></div> <div>66%</div> <div>14%</div> <div>•</div> <div>19%</div> </div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6873 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.

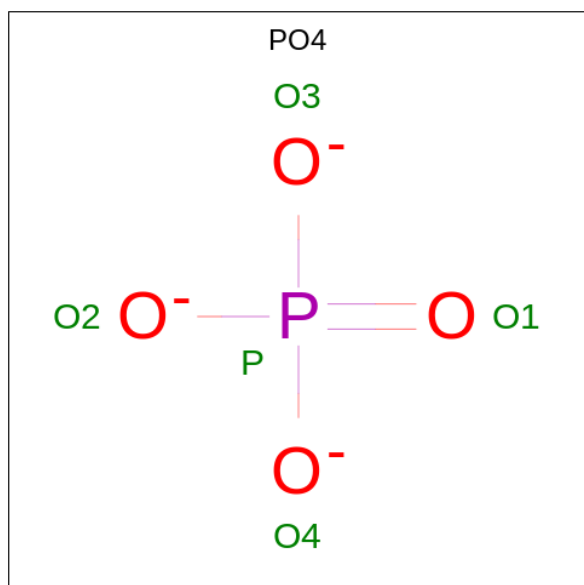
- Molecule 1 is a protein called Membrane-associated phosphatidylinositol transfer protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	0	0
			2345	1514	400	421	10			
1	B	303	Total	C	N	O	S	0	0	0
			2306	1488	393	416	9			
1	C	290	Total	C	N	O	S	0	0	0
			2216	1433	379	395	9			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	421	SER	-	expression tag	UNP O00562
B	421	SER	-	expression tag	UNP O00562
C	421	SER	-	expression tag	UNP O00562

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

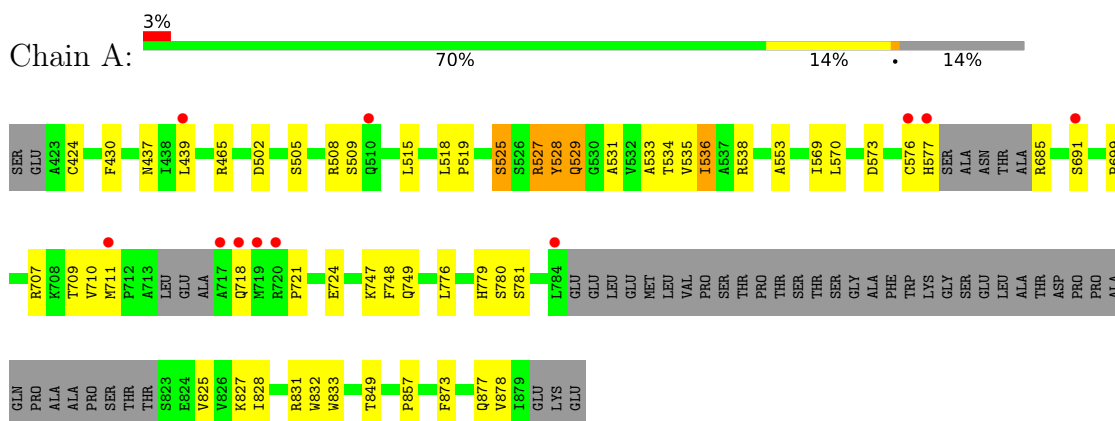
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		

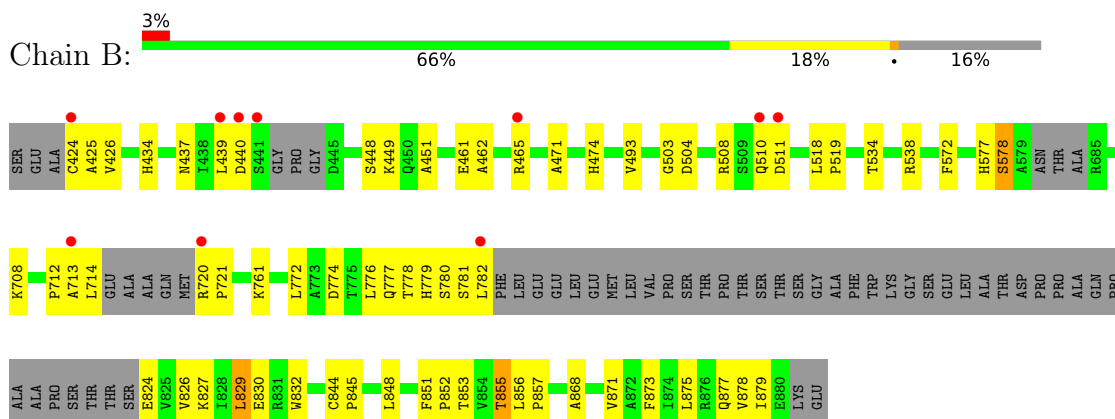
3 Residue-property plots

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.

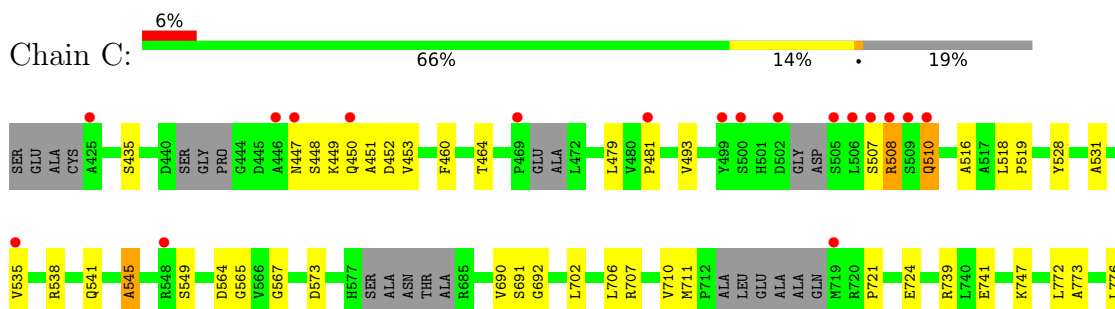
• Molecule 1: Membrane-associated phosphatidylinositol transfer protein 1

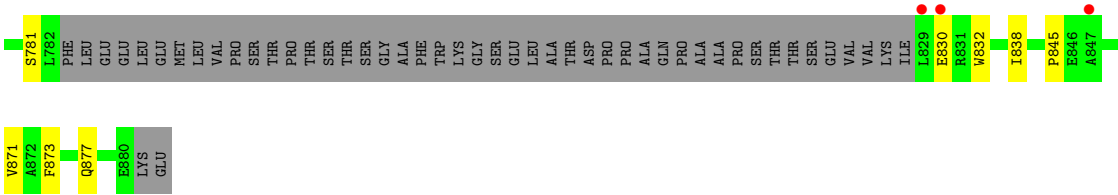


• Molecule 1: Membrane-associated phosphatidylinositol transfer protein 1



• Molecule 1: Membrane-associated phosphatidylinositol transfer protein 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	130.24Å 130.24Å 105.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.74 – 2.80 49.74 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.74-2.80) 97.2 (49.74-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.201 , 0.251 0.246 , 0.251	Depositor DCC
R_{free} test set	1993 reflections (7.70%)	wwPDB-VP
Wilson B-factor (Å ²)	76.4	Xtriage
Anisotropy	0.004	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 69.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6873	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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: PO4

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	0.46	1/2405 (0.0%)	0.91	2/3279 (0.1%)
1	B	0.41	0/2363	0.88	4/3221 (0.1%)
1	C	0.44	0/2271	0.93	4/3092 (0.1%)
All	All	0.44	1/7039 (0.0%)	0.91	10/9592 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	528	TYR	CA-C	-5.32	1.45	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	510	GLN	N-CA-C	-10.94	98.19	112.41
1	A	748	PHE	N-CA-C	-6.94	104.93	113.19
1	C	545	ALA	N-CA-C	-5.86	105.22	112.90
1	B	511	ASP	N-CA-C	5.70	118.50	110.23
1	A	709	THR	N-CA-C	5.61	117.48	111.36
1	B	829	LEU	N-CA-C	5.44	117.21	111.28
1	C	451	ALA	N-CA-C	-5.25	105.56	111.28
1	B	830	GLU	N-CA-C	5.20	117.35	111.11
1	C	845	PRO	CA-C-O	-5.07	115.56	121.34
1	B	578	SER	N-CA-C	5.03	117.17	110.53

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	2345	0	2344	41	0
1	B	2306	0	2305	55	0
1	C	2216	0	2213	48	0
2	A	5	0	0	0	0
3	A	1	0	0	0	0
All	All	6873	0	6862	136	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 (136) 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:845:PRO:HB2	1:C:510:GLN:HE22	1.14	1.09
1:A:849:THR:HB	1:B:510:GLN:HE21	1.35	0.91
1:C:450:GLN:HG2	1:C:479:LEU:HD22	1.52	0.89
1:C:450:GLN:HG2	1:C:479:LEU:CD2	2.04	0.86
1:A:439:LEU:HD11	1:A:857:PRO:HG2	1.57	0.86
1:B:845:PRO:HB2	1:C:510:GLN:NE2	1.90	0.86
1:B:713:ALA:O	1:B:714:LEU:HG	1.80	0.82
1:B:439:LEU:HD11	1:B:857:PRO:HG2	1.65	0.79
1:A:849:THR:HG21	1:B:510:GLN:HG2	1.65	0.79
1:A:849:THR:HB	1:B:510:GLN:NE2	1.99	0.78
1:B:848:LEU:HG	1:C:510:GLN:HE21	1.47	0.78
1:B:462:ALA:HA	1:B:465:ARG:HH21	1.49	0.77
1:C:447:ASN:C	1:C:449:LYS:H	1.95	0.74
1:C:453:VAL:HG21	1:C:479:LEU:HB2	1.69	0.74
1:C:481:PRO:HG2	1:C:538:ARG:HE	1.53	0.73
1:B:439:LEU:CD1	1:B:857:PRO:HG2	2.19	0.73
1:B:471:ALA:C	1:B:474:HIS:HD2	1.97	0.72
1:A:827:LYS:O	1:A:831:ARG:HG3	1.90	0.72
1:A:502:ASP:HB3	1:A:508:ARG:HD3	1.71	0.71
1:B:572:PHE:CE1	1:B:721:PRO:HG3	2.25	0.71
1:A:430:PHE:HZ	1:A:878:VAL:HG21	1.58	0.69
1:A:707:ARG:O	1:A:710:VAL:HG12	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:782:LEU:HD12	1:B:782:LEU:H	1.58	0.67
1:C:832:TRP:HE1	1:C:838:ILE:HG13	1.61	0.65
1:B:713:ALA:O	1:B:714:LEU:CG	2.45	0.64
1:B:824:GLU:O	1:B:827:LYS:HG2	1.98	0.64
1:C:450:GLN:HE21	1:C:479:LEU:HB3	1.61	0.64
1:B:824:GLU:O	1:B:827:LYS:NZ	2.31	0.63
1:A:533:ALA:HA	1:A:685:ARG:HH22	1.63	0.63
1:C:772:LEU:O	1:C:776:LEU:HG	1.99	0.62
1:C:447:ASN:C	1:C:449:LYS:N	2.56	0.62
1:B:471:ALA:C	1:B:474:HIS:CD2	2.76	0.62
1:A:528:TYR:O	1:A:529:GLN:C	2.41	0.62
1:A:691:SER:O	1:A:724:GLU:HB2	2.01	0.61
1:C:741:GLU:HG2	1:C:832:TRP:CZ2	2.35	0.61
1:A:849:THR:CG2	1:B:510:GLN:HG2	2.31	0.61
1:B:577:HIS:CE1	1:B:578:SER:O	2.55	0.60
1:B:782:LEU:HD12	1:B:782:LEU:N	2.18	0.58
1:A:439:LEU:HD11	1:A:857:PRO:CG	2.32	0.57
1:C:830:GLU:C	1:C:832:TRP:H	2.11	0.57
1:C:447:ASN:O	1:C:449:LYS:N	2.35	0.57
1:C:706:LEU:O	1:C:710:VAL:HG23	2.05	0.57
1:B:708:LYS:O	1:B:712:PRO:HA	2.05	0.56
1:C:747:LYS:NZ	1:C:781:SER:O	2.22	0.56
1:C:460:PHE:O	1:C:464:THR:HG23	2.06	0.56
1:A:779:HIS:O	1:A:781:SER:N	2.38	0.55
1:C:450:GLN:NE2	1:C:479:LEU:HB3	2.22	0.55
1:A:439:LEU:CD1	1:A:857:PRO:HG2	2.31	0.54
1:B:826:VAL:HA	1:B:829:LEU:HD23	1.89	0.54
1:C:691:SER:O	1:C:724:GLU:HB2	2.08	0.54
1:A:873:PHE:O	1:A:877:GLN:HG2	2.09	0.53
1:B:772:LEU:O	1:B:776:LEU:HG	2.08	0.52
1:C:707:ARG:O	1:C:711:MET:N	2.33	0.52
1:A:710:VAL:HG13	1:A:711:MET:N	2.24	0.52
1:A:779:HIS:O	1:A:780:SER:C	2.53	0.52
1:A:430:PHE:CZ	1:A:878:VAL:HG21	2.43	0.51
1:B:774:ASP:O	1:B:778:THR:HG23	2.10	0.51
1:B:848:LEU:HG	1:C:510:GLN:NE2	2.20	0.51
1:C:435:SER:N	1:C:564:ASP:OD2	2.43	0.51
1:B:503:GLY:O	1:B:508:ARG:NH2	2.43	0.51
1:A:528:TYR:O	1:A:531:ALA:N	2.44	0.51
1:C:773:ALA:HA	1:C:776:LEU:HD12	1.93	0.51
1:C:739:ARG:NH2	1:C:741:GLU:OE2	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:434:HIS:O	1:B:449:LYS:NZ	2.36	0.50
1:B:461:GLU:O	1:B:465:ARG:HG3	2.11	0.50
1:C:873:PHE:O	1:C:877:GLN:HG2	2.11	0.50
1:B:471:ALA:O	1:B:474:HIS:HB2	2.12	0.49
1:B:781:SER:O	1:B:782:LEU:C	2.54	0.49
1:C:528:TYR:OH	1:C:573:ASP:OD2	2.26	0.49
1:B:777:GLN:O	1:B:780:SER:OG	2.27	0.49
1:B:448:SER:O	1:B:451:ALA:HB3	2.12	0.49
1:B:851:PHE:HB2	1:B:856:LEU:HD21	1.95	0.48
1:C:832:TRP:NE1	1:C:838:ILE:HG13	2.27	0.48
1:A:828:ILE:O	1:A:832:TRP:HB2	2.13	0.48
1:A:465:ARG:O	1:A:465:ARG:HD3	2.14	0.47
1:A:776:LEU:HD13	1:A:825:VAL:HG13	1.97	0.47
1:B:462:ALA:HA	1:B:465:ARG:NH2	2.25	0.47
1:B:572:PHE:HE1	1:B:721:PRO:HG3	1.75	0.47
1:B:761:LYS:HB3	1:B:761:LYS:HE3	1.67	0.47
1:B:875:LEU:HA	1:B:878:VAL:HG22	1.96	0.47
1:B:873:PHE:O	1:B:877:GLN:HG2	2.15	0.47
1:C:690:VAL:CG1	1:C:692:GLY:O	2.63	0.47
1:B:844:CYS:SG	1:B:848:LEU:HD12	2.55	0.46
1:A:502:ASP:HB3	1:A:508:ARG:CD	2.44	0.46
1:B:462:ALA:CA	1:B:465:ARG:HH21	2.26	0.46
1:A:424:CYS:HB2	1:A:553:ALA:O	2.16	0.46
1:C:450:GLN:CG	1:C:479:LEU:HD22	2.35	0.46
1:C:531:ALA:O	1:C:535:VAL:HG23	2.16	0.46
1:A:439:LEU:HD21	1:A:515:LEU:HD21	1.98	0.46
1:B:852:PRO:O	1:B:855:THR:HG22	2.15	0.46
1:B:776:LEU:O	1:B:779:HIS:O	2.33	0.46
1:B:782:LEU:N	1:B:782:LEU:CD1	2.79	0.45
1:C:516:ALA:O	1:C:519:PRO:HD2	2.15	0.45
1:A:535:VAL:HG21	1:A:570:LEU:HB3	1.97	0.45
1:C:741:GLU:HG2	1:C:832:TRP:HZ2	1.81	0.45
1:A:505:SER:O	1:A:509:SER:HB3	2.16	0.45
1:C:724:GLU:N	1:C:724:GLU:OE1	2.51	0.44
1:C:564:ASP:OD1	1:C:567:GLY:N	2.47	0.44
1:B:504:ASP:O	1:B:508:ARG:HG3	2.18	0.44
1:A:747:LYS:C	1:A:749:GLN:N	2.76	0.44
1:C:564:ASP:O	1:C:565:GLY:C	2.61	0.44
1:A:747:LYS:HA	1:A:747:LYS:HD3	1.71	0.44
1:B:868:ALA:HA	1:B:871:VAL:HG22	2.00	0.43
1:C:545:ALA:O	1:C:549:SER:HB3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:830:GLU:C	1:C:832:TRP:N	2.75	0.43
1:B:503:GLY:HA2	1:B:508:ARG:HH21	1.83	0.43
1:A:569:ILE:HD11	1:A:699:PRO:HG2	2.01	0.43
1:A:576:CYS:O	1:A:577:HIS:HB3	2.19	0.42
1:C:507:SER:O	1:C:508:ARG:C	2.61	0.42
1:B:875:LEU:O	1:B:879:ILE:HG23	2.19	0.42
1:C:460:PHE:HA	1:C:871:VAL:HG11	2.01	0.42
1:A:437:ASN:HD21	1:A:439:LEU:HB2	1.84	0.42
1:A:534:THR:O	1:A:538:ARG:HG3	2.18	0.42
1:B:437:ASN:HB3	1:B:440:ASP:OD1	2.20	0.42
1:C:481:PRO:HG2	1:C:538:ARG:NE	2.27	0.42
1:A:721:PRO:HD2	1:A:833:TRP:CD1	2.54	0.42
1:B:772:LEU:HD22	1:B:832:TRP:CG	2.54	0.42
1:A:529:GLN:H	1:A:529:GLN:HG3	1.48	0.42
1:C:450:GLN:HG2	1:C:479:LEU:HD23	1.95	0.42
1:B:534:THR:O	1:B:538:ARG:HG3	2.20	0.42
1:B:424:CYS:O	1:B:425:ALA:HB2	2.20	0.42
1:A:779:HIS:C	1:A:781:SER:N	2.73	0.42
1:C:518:LEU:HB2	1:C:519:PRO:HD3	2.02	0.41
1:C:538:ARG:HA	1:C:541:GLN:HB3	2.02	0.41
1:A:525:SER:OG	1:A:527:ARG:HG2	2.20	0.41
1:C:449:LYS:O	1:C:452:ASP:HB2	2.19	0.41
1:C:516:ALA:HB1	1:C:702:LEU:CD2	2.50	0.41
1:C:773:ALA:O	1:C:776:LEU:HB2	2.21	0.41
1:A:518:LEU:HB2	1:A:519:PRO:HD3	2.02	0.41
1:B:720:ARG:HH11	1:B:720:ARG:HG3	1.86	0.41
1:A:528:TYR:OH	1:A:573:ASP:OD2	2.36	0.40
1:A:536:ILE:HD12	1:A:685:ARG:NE	2.36	0.40
1:C:450:GLN:CG	1:C:479:LEU:CD2	2.89	0.40
1:B:518:LEU:HB2	1:B:519:PRO:HD3	2.03	0.40
1:B:772:LEU:HD22	1:B:832:TRP:CD2	2.57	0.40
1:B:824:GLU:HA	1:B:824:GLU:OE1	2.22	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	301/360 (84%)	294 (98%)	7 (2%)	0	100	100
1	B	293/360 (81%)	289 (99%)	4 (1%)	0	100	100
1	C	276/360 (77%)	267 (97%)	7 (2%)	2 (1%)	19	48
All	All	870/1080 (81%)	850 (98%)	18 (2%)	2 (0%)	44	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	448	SER
1	C	721	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	249/290 (86%)	244 (98%)	5 (2%)	50	81
1	B	246/290 (85%)	242 (98%)	4 (2%)	58	85
1	C	236/290 (81%)	234 (99%)	2 (1%)	79	93
All	All	731/870 (84%)	720 (98%)	11 (2%)	60	86

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	525	SER
1	A	527	ARG
1	A	529	GLN
1	A	536	ILE
1	A	718	GLN
1	B	426	VAL
1	B	493	VAL

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Mol	Chain	Res	Type
1	B	853	THR
1	B	855	THR
1	C	493	VAL
1	C	508	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	779	HIS
1	B	474	HIS
1	B	510	GLN
1	B	577	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 ⓘ

1 ligand is 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	$\# Z > 2$
2	PO4	A	901	-	4,4,4	0.91	0	6,6,6	0.45	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	309/360 (85%)	0.06	11 (3%)	46 38	52, 75, 123, 195	0
1	B	303/360 (84%)	0.14	10 (3%)	49 41	50, 85, 135, 166	0
1	C	290/360 (80%)	0.54	21 (7%)	23 17	56, 105, 159, 221	0
All	All	902/1080 (83%)	0.24	42 (4%)	37 30	50, 87, 149, 221	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	719	MET	5.0
1	A	717	ALA	4.9
1	A	720	ARG	4.8
1	A	577	HIS	4.7
1	A	711	MET	4.6
1	B	782	LEU	4.6
1	C	502	ASP	4.0
1	C	719	MET	4.0
1	C	829	LEU	3.9
1	C	505	SER	3.9
1	C	510	GLN	3.6
1	B	465	ARG	3.4
1	B	510	GLN	3.4
1	C	508	ARG	3.4
1	A	718	GLN	3.1
1	B	440	ASP	2.9
1	C	507	SER	2.9
1	C	500	SER	2.8
1	C	509	SER	2.8
1	C	450	GLN	2.8
1	C	499	TYR	2.7
1	A	576	CYS	2.7
1	C	548	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	511	ASP	2.7
1	B	424	CYS	2.7
1	C	469	PRO	2.6
1	C	446	ALA	2.6
1	C	830	GLU	2.5
1	B	441	SER	2.5
1	C	535	VAL	2.5
1	A	691	SER	2.4
1	A	784	LEU	2.4
1	B	439	LEU	2.4
1	C	447	ASN	2.4
1	C	481	PRO	2.3
1	C	425	ALA	2.3
1	C	506	LEU	2.3
1	A	510	GLN	2.1
1	B	720	ARG	2.1
1	A	439	LEU	2.0
1	B	713	ALA	2.0
1	C	847	ALA	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(\AA^2)	Q<0.9
2	PO4	A	901	5/5	0.77	0.10	120,125,126,130	0

6.5 Other polymers [i](#)

There are no such residues in this entry.