



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:32 AM UTC

PDB ID : 1QAS / pdb_00001qas
Title : 1-PHOSPHATIDYLINOSITOL-4,5-BISPHOSPHATE PHOSPHODI-
ESTERASE DELTA 1
Authors : Grobler, J.A.; Hurley, J.H.
Deposited on : 1996-08-02
Resolution : 2.40 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
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

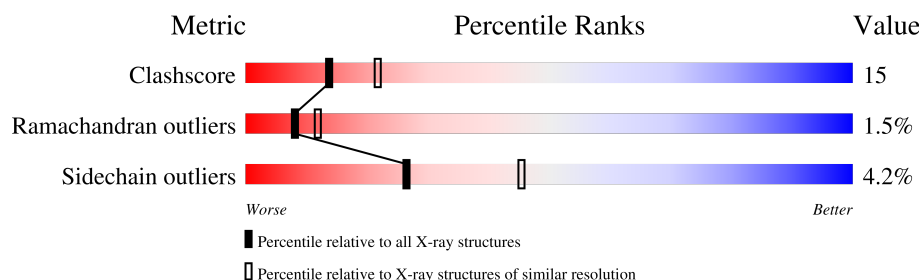
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	622	
1	B	622	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8235 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 PHOSPHOLIPASE C DELTA-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	505	Total	C	N	O	S	0	0	0
			3990	2522	696	750	22			
1	B	504	Total	C	N	O	S	0	0	0
			3979	2517	696	744	22			

- Molecule 2 is water.

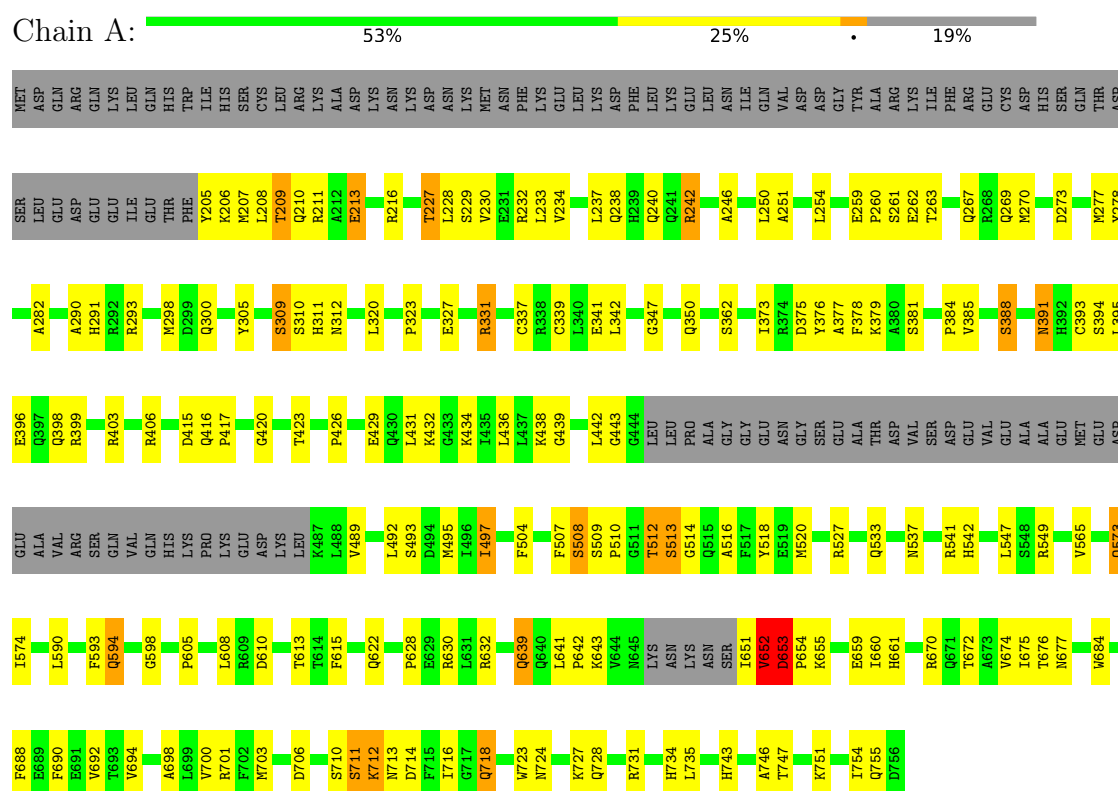
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	132	Total	O	0	0
			132	132		
2	B	134	Total	O	0	0
			134	134		

3 Residue-property plots

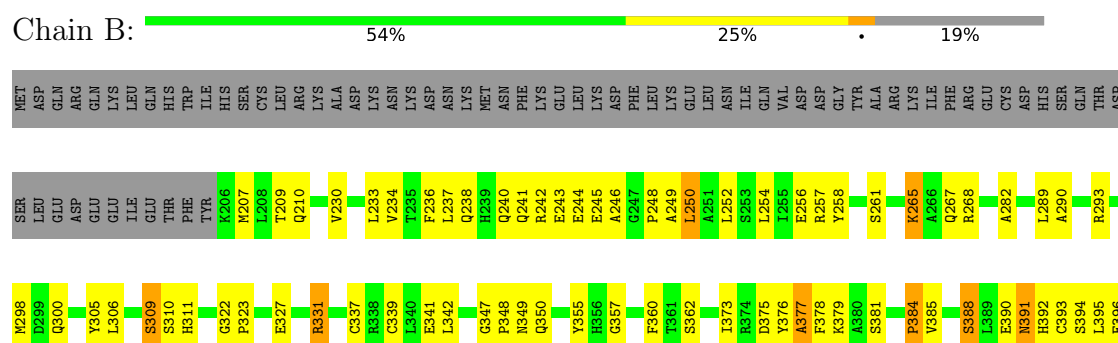
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

• Molecule 1: PHOSPHOLIPASE C DELTA-1



• Molecule 1: PHOSPHOLIPASE C DELTA-1



Q397	Q398	R399	R403	D415	Q416	T423	P426	E429	Q430	L431	K432	Q433	K434	L435	L436	L437	K438	Q439	G443	G444	L445	LEU	PRO	ALA	GLY	GLY	GLU	ASN	GLY	SER	GLU	ALA	THR	ASP	ASP	VAL	SER	SER	ASP	GLU	VAL	GLU	ALA	ALA	GLU	GLU	MET	GLU	ASP	GLU	ALA	VAL	ARG	SER	GLN	VAL
GLN	HIS	LYS	PRO	LYS	GLU	ASP	K485	V489	L492	M495	I496	I497	K500	F504	F507	S508	S509	P510	GLY	THR	SER	G514	Q515	A516	F517	Y518	E519	M520	S524	E525	S526	R527	N537	G538	L547	S548	R549	P552	A553	V565	Q573	I574	L590	F593	Q594											
G598	P605	L608	R609	D610	T613	T614	F615	R616	S617	Q622	P628	R629	R630	I636	Q639	G640	L641	P642	K643	V644	N645	LYS	ASN	LYS	ASN	SER	I651	V652	D653	P654	E659	I660	H661	G662	R670	Q671	T672	I675	T676	N677	W684	D685	M686	E687	F688	E689	F690									
E691	V692	T693	V694	A698	L699	V700	R701	M703	D706	S710	S711	K712	N713	D714	Q718	W723	K727	Q728	G729	Y730	R731	H734	L735	H743	A746	I754	Q755	D756																												

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	61.10Å 75.40Å 86.90Å 66.90° 85.40° 89.80°	Depositor
Resolution (Å)	6.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.40)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.212 , 0.286	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8235	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	0.63	0/4083	1.07	21/5531 (0.4%)
1	B	0.63	1/4070 (0.0%)	1.09	28/5511 (0.5%)
All	All	0.63	1/8153 (0.0%)	1.08	49/11042 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	686	MET	SD-CE	5.21	1.92	1.79

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	244	GLU	N-CA-C	-11.27	95.58	111.81
1	A	514	GLY	N-CA-C	8.99	122.64	110.69
1	B	594	GLN	N-CA-C	-8.57	101.44	112.23
1	A	594	GLN	N-CA-C	-7.35	102.97	112.23
1	A	391	ASN	N-CA-C	7.21	120.28	109.25
1	B	258	TYR	N-CA-C	7.15	122.09	113.23
1	A	331	ARG	NE-CZ-NH2	-7.11	112.80	119.20
1	B	384	PRO	N-CA-C	6.88	122.02	111.15
1	A	443	GLY	N-CA-C	-6.84	96.98	113.18
1	A	331	ARG	CD-NE-CZ	6.80	133.93	124.40
1	A	384	PRO	N-CA-C	6.77	121.85	111.15
1	B	391	ASN	N-CA-C	6.58	119.32	109.25
1	A	309	SER	N-CA-C	6.46	119.68	108.76
1	B	331	ARG	NE-CZ-NH2	-6.22	113.60	119.20
1	A	497	ILE	N-CA-C	6.20	116.89	111.56
1	A	331	ARG	NE-CZ-NH1	6.14	127.64	121.50
1	B	309	SER	N-CA-C	6.12	119.10	108.76
1	B	331	ARG	CD-NE-CZ	6.11	132.95	124.40
1	B	489	VAL	CA-C-N	6.08	126.25	119.32
1	B	489	VAL	C-N-CA	6.08	126.25	119.32
1	A	694	VAL	CA-C-N	6.00	125.88	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	694	VAL	C-N-CA	6.00	125.88	119.28
1	B	265	LYS	N-CA-C	-5.89	104.78	111.14
1	B	694	VAL	CA-C-N	5.87	125.73	119.28
1	B	694	VAL	C-N-CA	5.87	125.73	119.28
1	B	654	PRO	N-CA-C	5.76	120.33	111.57
1	B	375	ASP	N-CA-C	5.74	117.62	111.36
1	B	565	VAL	N-CA-C	5.50	115.95	110.23
1	A	565	VAL	N-CA-C	5.46	115.91	110.23
1	B	250	LEU	N-CA-C	-5.45	105.24	111.07
1	B	641	LEU	CA-C-N	5.43	125.19	119.76
1	B	641	LEU	C-N-CA	5.43	125.19	119.76
1	A	654	PRO	N-CA-C	5.32	119.66	111.57
1	B	306	LEU	N-CA-C	-5.30	100.60	109.24
1	A	653	ASP	N-CA-C	5.26	121.44	109.81
1	B	538	GLY	N-CA-C	-5.26	106.03	112.77
1	A	320	LEU	N-CA-C	5.26	119.92	112.93
1	B	257	ARG	N-CA-C	5.26	116.86	111.03
1	A	735	LEU	N-CA-C	5.22	117.45	110.35
1	A	375	ASP	N-CA-C	5.21	117.03	111.36
1	A	282	ALA	N-CA-C	-5.18	105.72	111.36
1	A	347	GLY	CA-C-N	5.12	125.41	119.93
1	A	347	GLY	C-N-CA	5.12	125.41	119.93
1	B	735	LEU	N-CA-C	5.11	117.71	110.50
1	B	347	GLY	CA-C-N	5.11	125.40	119.93
1	B	347	GLY	C-N-CA	5.11	125.40	119.93
1	B	524	SER	N-CA-C	-5.08	103.69	110.55
1	B	377	ALA	N-CA-C	5.06	117.45	111.33
1	B	497	ILE	N-CA-C	5.02	116.20	111.48

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	3990	0	3906	123	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3979	0	3903	111	0
2	A	132	0	0	5	0
2	B	134	0	0	7	0
All	All	8235	0	7809	229	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 (229) 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:228:LEU:HB2	1:A:270:MET:HB3	1.47	0.94
1:A:520:MET:HE3	1:A:549:ARG:HB2	1.46	0.94
1:B:520:MET:HE3	1:B:549:ARG:HB2	1.50	0.92
1:B:438:LYS:HE3	1:B:520:MET:HE2	1.58	0.84
1:A:399:ARG:O	1:A:403:ARG:HG2	1.79	0.82
1:A:573:GLN:H	1:A:573:GLN:HE21	1.31	0.79
1:B:399:ARG:O	1:B:403:ARG:HG2	1.82	0.79
1:B:241:GLN:HE22	1:B:730:TYR:H	1.31	0.77
1:B:734:HIS:HE1	2:B:2534:HOH:O	1.68	0.76
1:B:282:ALA:HB1	1:B:289:LEU:HD13	1.66	0.76
1:B:207:MET:HA	1:B:210:GLN:NE2	2.00	0.75
1:A:590:LEU:O	1:A:594:GLN:HG2	1.87	0.74
1:A:547:LEU:HD23	1:A:573:GLN:HG3	1.70	0.74
1:B:246:ALA:HB3	1:B:250:LEU:HB2	1.71	0.73
1:B:339:CYS:HB3	2:B:2946:HOH:O	1.89	0.72
1:B:323:PRO:HA	1:B:362:SER:HB3	1.73	0.70
1:A:701:ARG:HE	1:A:718:GLN:HE21	1.38	0.69
1:A:227:THR:CG2	1:A:269:GLN:HB3	2.23	0.69
1:B:520:MET:HE3	1:B:549:ARG:CB	2.22	0.69
1:B:701:ARG:HE	1:B:718:GLN:HE21	1.38	0.69
1:B:391:ASN:HD21	1:B:393:CYS:HB2	1.58	0.68
1:A:520:MET:HE3	1:A:549:ARG:CB	2.22	0.68
1:B:309:SER:OG	1:B:574:ILE:HG23	1.93	0.68
1:B:507:PHE:O	1:B:508:SER:HB2	1.94	0.68
1:B:573:GLN:H	1:B:573:GLN:HE21	1.42	0.68
1:A:339:CYS:HB3	2:A:2235:HOH:O	1.94	0.68
1:B:547:LEU:HD23	1:B:573:GLN:HG3	1.76	0.67
1:A:439:GLY:HA2	2:A:2318:HOH:O	1.94	0.67
1:B:298:MET:HB2	1:B:429:GLU:HG2	1.76	0.66
1:B:590:LEU:O	1:B:594:GLN:HG2	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:THR:HG21	1:A:269:GLN:HB3	1.78	0.66
1:A:395:LEU:HD22	1:A:489:VAL:HG13	1.77	0.66
1:A:573:GLN:H	1:A:573:GLN:NE2	1.93	0.65
1:A:298:MET:HB2	1:A:429:GLU:HG2	1.77	0.65
1:B:617:SER:HB3	2:B:2890:HOH:O	1.97	0.65
1:A:642:PRO:HG3	1:A:743:HIS:CE1	2.32	0.65
1:A:438:LYS:HE3	1:A:520:MET:HE2	1.80	0.64
1:B:243:GLU:C	1:B:245:GLU:H	2.05	0.64
1:A:323:PRO:HA	1:A:362:SER:HB3	1.79	0.64
1:A:327:GLU:O	1:A:331:ARG:HG3	1.98	0.64
1:A:309:SER:OG	1:A:574:ILE:HG23	1.98	0.63
1:B:429:GLU:O	1:B:432:LYS:HG3	1.98	0.62
1:B:672:THR:HG22	1:B:688:PHE:HZ	1.63	0.62
1:A:420:GLY:H	1:B:349:ASN:ND2	1.98	0.62
1:A:642:PRO:HD3	1:A:746:ALA:CB	2.30	0.61
1:B:350:GLN:NE2	1:B:396:GLU:HG3	2.16	0.61
1:B:651:ILE:HG22	1:B:652:VAL:H	1.65	0.60
1:A:230:VAL:O	1:A:234:VAL:HG23	2.02	0.60
1:A:512:THR:O	1:A:513:SER:HB2	2.02	0.60
1:B:246:ALA:O	1:B:249:ALA:HB3	2.02	0.60
1:A:417:PRO:HG2	1:B:348:PRO:HB2	1.84	0.59
1:A:395:LEU:O	1:A:399:ARG:HG3	2.02	0.59
1:A:643:LYS:CB	1:A:651:ILE:HB	2.33	0.59
1:A:260:PRO:HG3	1:A:273:ASP:HB3	1.85	0.59
1:A:350:GLN:NE2	1:A:396:GLU:HG3	2.18	0.59
1:B:444:GLY:H	1:B:500:LYS:NZ	2.02	0.58
1:A:520:MET:CE	1:A:549:ARG:HB2	2.27	0.58
1:B:573:GLN:H	1:B:573:GLN:NE2	2.01	0.58
1:B:230:VAL:HG23	1:B:268:ARG:O	2.04	0.58
1:B:395:LEU:O	1:B:399:ARG:HG3	2.04	0.58
1:B:672:THR:HG22	1:B:688:PHE:CZ	2.39	0.58
1:A:237:LEU:HD13	1:A:250:LEU:HG	1.86	0.57
1:A:238:GLN:HB3	1:A:246:ALA:CB	2.34	0.57
1:B:350:GLN:HE22	1:B:396:GLU:HG3	1.69	0.57
1:B:727:LYS:HE3	1:B:731:ARG:NH2	2.20	0.57
1:A:643:LYS:C	1:A:651:ILE:HG13	2.29	0.57
1:A:630:ARG:NH1	1:A:630:ARG:HB2	2.19	0.57
1:B:537:ASN:HB3	1:B:615:PHE:O	2.05	0.57
1:A:672:THR:HG22	1:A:688:PHE:HZ	1.69	0.57
1:B:426:PRO:HG2	1:B:431:LEU:HD11	1.86	0.57
1:B:256:GLU:O	1:B:265:LYS:HE3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:755:GLN:HG2	2:B:2656:HOH:O	2.06	0.56
1:A:605:PRO:HD2	1:A:608:LEU:HD12	1.85	0.56
1:A:701:ARG:HE	1:A:718:GLN:NE2	2.03	0.55
1:B:327:GLU:O	1:B:331:ARG:HG3	2.06	0.55
1:B:652:VAL:O	1:B:654:PRO:HD3	2.05	0.55
1:A:642:PRO:HD3	1:A:746:ALA:HB2	1.87	0.55
1:B:394:SER:O	1:B:398:GLN:HG3	2.06	0.55
1:A:533:GLN:NE2	2:A:2320:HOH:O	2.40	0.55
1:A:426:PRO:HG2	1:A:431:LEU:HD11	1.89	0.55
1:B:605:PRO:HD2	1:B:608:LEU:HD12	1.88	0.55
1:A:504:PHE:HB3	1:A:527:ARG:HH22	1.72	0.54
1:A:350:GLN:HE22	1:A:396:GLU:HG3	1.72	0.54
1:A:672:THR:HG22	1:A:688:PHE:CZ	2.43	0.54
1:A:675:ILE:HG12	1:A:684:TRP:NE1	2.22	0.54
1:B:243:GLU:C	1:B:245:GLU:N	2.65	0.54
1:B:241:GLN:HE22	1:B:730:TYR:N	2.04	0.54
1:B:520:MET:CE	1:B:549:ARG:HB2	2.33	0.54
1:A:710:SER:O	1:A:711:SER:HB2	2.08	0.54
1:B:234:VAL:O	1:B:238:GLN:HG2	2.08	0.54
1:A:391:ASN:HD21	1:A:393:CYS:HB2	1.72	0.54
1:A:211:ARG:HB3	1:A:213:GLU:OE2	2.07	0.53
1:B:261:SER:O	1:B:265:LYS:HB2	2.08	0.53
1:A:406:ARG:NH2	1:B:360:PHE:O	2.42	0.53
1:A:547:LEU:CD2	1:A:573:GLN:HG3	2.36	0.53
1:B:651:ILE:HG22	1:B:652:VAL:N	2.23	0.53
1:A:728:GLN:NE2	1:A:754:ILE:H	2.06	0.53
1:B:230:VAL:O	1:B:234:VAL:HG23	2.07	0.53
1:B:659:GLU:HG3	1:B:701:ARG:HB3	1.90	0.53
1:A:727:LYS:HE3	1:A:731:ARG:NH2	2.24	0.53
1:A:298:MET:CB	1:A:429:GLU:HG2	2.40	0.52
1:B:298:MET:CB	1:B:429:GLU:HG2	2.40	0.52
1:B:376:TYR:HA	1:B:379:LYS:HG2	1.90	0.52
1:B:710:SER:O	1:B:711:SER:HB2	2.09	0.52
1:A:642:PRO:HG2	1:A:716:ILE:CG2	2.40	0.52
1:B:675:ILE:HG12	1:B:684:TRP:NE1	2.24	0.52
1:B:630:ARG:NH1	1:B:630:ARG:HB2	2.25	0.52
1:A:228:LEU:HD13	1:A:233:LEU:HD13	1.92	0.52
1:A:659:GLU:HG3	1:A:701:ARG:HB3	1.90	0.52
1:A:376:TYR:HA	1:A:379:LYS:HG2	1.90	0.52
1:A:234:VAL:O	1:A:238:GLN:HG2	2.10	0.51
1:A:240:GLN:HE22	1:A:751:LYS:NZ	2.07	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:388:SER:HA	1:A:438:LYS:HB3	1.92	0.51
1:B:701:ARG:HE	1:B:718:GLN:NE2	2.05	0.51
1:A:610:ASP:O	1:A:613:THR:HG22	2.10	0.51
1:A:670:ARG:HG2	1:A:690:PHE:CE2	2.46	0.51
1:B:311:HIS:CD2	1:B:552:PRO:HD2	2.45	0.51
1:A:420:GLY:N	1:B:349:ASN:HD21	2.09	0.51
1:B:415:ASP:C	1:B:497:ILE:HD11	2.36	0.51
1:B:728:GLN:NE2	1:B:754:ILE:H	2.08	0.51
1:A:442:LEU:HB2	1:A:493:SER:OG	2.11	0.50
1:A:227:THR:HG22	1:A:270:MET:N	2.26	0.50
1:B:388:SER:HA	1:B:438:LYS:HB3	1.93	0.50
1:A:415:ASP:O	1:A:497:ILE:HD11	2.12	0.50
1:A:238:GLN:HB3	1:A:246:ALA:HB3	1.93	0.50
1:A:420:GLY:H	1:B:349:ASN:HD21	1.59	0.50
1:B:547:LEU:CD2	1:B:573:GLN:HG3	2.40	0.50
1:B:662:GLY:HA3	2:B:2796:HOH:O	2.11	0.49
1:B:652:VAL:HG12	1:B:654:PRO:HG3	1.94	0.49
1:B:246:ALA:H	1:B:250:LEU:HB2	1.77	0.49
1:A:240:GLN:HE22	1:A:751:LYS:HZ1	1.60	0.48
1:B:672:THR:CG2	1:B:688:PHE:HZ	2.27	0.48
1:A:290:ALA:O	1:A:293:ARG:HG2	2.13	0.48
1:A:429:GLU:O	1:A:432:LYS:HG3	2.12	0.48
1:B:439:GLY:HA2	2:B:3008:HOH:O	2.13	0.48
1:B:610:ASP:O	1:B:613:THR:HG22	2.14	0.48
1:A:692:VAL:HG21	1:A:723:TRP:CZ3	2.48	0.48
1:B:516:ALA:HB1	1:B:518:TYR:CE1	2.49	0.48
1:B:373:ILE:O	1:B:377:ALA:HB2	2.14	0.47
1:A:228:LEU:O	1:A:269:GLN:HA	2.13	0.47
1:A:507:PHE:O	1:A:508:SER:HB3	2.13	0.47
1:A:628:PRO:HA	1:A:692:VAL:O	2.15	0.47
1:A:639:GLN:HG2	1:A:747:THR:OG1	2.15	0.47
1:A:661:HIS:O	1:A:698:ALA:HA	2.14	0.47
1:B:350:GLN:HE22	1:B:396:GLU:CG	2.28	0.47
1:A:573:GLN:NE2	1:A:573:GLN:N	2.62	0.46
1:B:310:SER:HB2	1:B:337:CYS:SG	2.55	0.46
1:B:438:LYS:HE3	1:B:520:MET:CE	2.37	0.46
1:A:205:TYR:HB3	1:A:209:THR:OG1	2.15	0.46
1:A:350:GLN:HE22	1:A:396:GLU:CG	2.29	0.46
1:A:651:ILE:O	1:A:652:VAL:HG13	2.16	0.46
1:A:652:VAL:HG11	1:A:677:ASN:HD22	1.80	0.46
1:A:492:LEU:HA	1:A:495:MET:HE2	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:661:HIS:O	1:B:698:ALA:HA	2.15	0.46
1:A:227:THR:HB	1:A:269:GLN:NE2	2.30	0.46
1:A:632:ARG:HH12	1:A:755:GLN:HG3	1.80	0.46
1:B:300:GLN:HB2	1:B:305:TYR:CE1	2.50	0.46
1:B:236:PHE:CD1	1:B:240:GLN:HG3	2.50	0.45
1:B:692:VAL:HG21	1:B:723:TRP:CZ3	2.51	0.45
1:A:254:LEU:HD22	1:A:278:TYR:CD1	2.52	0.45
1:A:373:ILE:O	1:A:377:ALA:HB2	2.17	0.45
1:B:642:PRO:HD3	1:B:746:ALA:HB2	1.99	0.45
1:B:712:LYS:HD3	1:B:713:ASN:N	2.32	0.45
1:A:712:LYS:HD3	1:A:713:ASN:N	2.32	0.45
1:B:290:ALA:O	1:B:293:ARG:HG2	2.17	0.45
1:B:670:ARG:HG2	1:B:690:PHE:CE2	2.51	0.45
1:A:341:GLU:C	1:A:342:LEU:HD12	2.42	0.44
1:A:672:THR:CG2	1:A:688:PHE:HZ	2.30	0.44
1:B:415:ASP:O	1:B:497:ILE:HD11	2.17	0.44
1:B:492:LEU:HA	1:B:495:MET:HE2	1.98	0.44
1:A:205:TYR:HA	1:A:208:LEU:HB3	1.99	0.44
1:B:660:ILE:CD1	1:B:700:VAL:HG22	2.48	0.44
1:B:378:PHE:HA	1:B:381:SER:O	2.18	0.44
1:B:504:PHE:HB3	1:B:527:ARG:HH22	1.82	0.44
1:A:263:THR:O	1:A:267:GLN:HG3	2.18	0.44
1:B:628:PRO:HA	1:B:692:VAL:O	2.18	0.44
1:A:642:PRO:HG2	1:A:716:ILE:HG23	1.99	0.43
1:B:384:PRO:HG3	1:B:431:LEU:HB2	2.00	0.43
1:A:537:ASN:HB3	1:A:615:PHE:O	2.18	0.43
1:A:259:GLU:OE2	1:A:261:SER:HB3	2.18	0.43
1:A:652:VAL:HG11	1:A:677:ASN:ND2	2.33	0.43
1:B:525:GLU:HG2	1:B:553:ALA:HB2	2.00	0.43
1:A:310:SER:HB2	1:A:337:CYS:SG	2.59	0.43
1:A:734:HIS:HE1	2:A:1821:HOH:O	2.01	0.43
1:B:341:GLU:C	1:B:342:LEU:HD12	2.44	0.43
1:A:438:LYS:HE3	1:A:520:MET:CE	2.47	0.43
1:B:355:TYR:CZ	1:B:357:GLY:HA2	2.54	0.42
1:A:415:ASP:C	1:A:497:ILE:HD11	2.43	0.42
1:B:594:GLN:HA	1:B:594:GLN:NE2	2.34	0.42
1:A:537:ASN:O	1:A:541:ARG:HG3	2.19	0.42
1:A:593:PHE:O	1:A:598:GLY:HA2	2.19	0.42
1:A:311:HIS:O	1:A:312:ASN:C	2.62	0.42
1:A:395:LEU:HD22	1:A:489:VAL:CG1	2.47	0.42
1:A:504:PHE:HB3	1:A:527:ARG:NH2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:241:GLN:HA	1:B:241:GLN:NE2	2.35	0.42
1:A:655:LYS:HD2	1:A:674:VAL:HG22	2.00	0.42
1:A:706:ASP:HB3	1:A:714:ASP:HB2	2.00	0.42
1:A:238:GLN:O	1:A:242:ARG:HA	2.19	0.42
1:B:322:GLY:HA3	2:B:2504:HOH:O	2.19	0.42
1:A:728:GLN:HE21	1:A:754:ILE:H	1.68	0.42
1:A:660:ILE:CD1	1:A:700:VAL:HG22	2.50	0.42
1:B:653:ASP:OD2	1:B:676:THR:HA	2.19	0.42
1:B:207:MET:HA	1:B:210:GLN:HE21	1.83	0.42
1:A:206:LYS:HA	1:A:210:GLN:HG3	2.01	0.41
1:A:233:LEU:HD23	1:A:251:ALA:HB1	2.02	0.41
1:A:653:ASP:OD2	1:A:676:THR:HA	2.19	0.41
1:B:248:PRO:O	1:B:252:LEU:HG	2.20	0.41
1:A:229:SER:O	1:A:230:VAL:C	2.63	0.41
1:A:516:ALA:HB1	1:A:518:TYR:CE2	2.56	0.41
1:A:542:HIS:HE1	2:A:2411:HOH:O	2.02	0.41
1:B:641:LEU:HD12	1:B:641:LEU:HA	1.88	0.41
1:B:706:ASP:HB3	1:B:714:ASP:HB2	2.01	0.41
1:A:632:ARG:NH1	1:A:755:GLN:HG3	2.35	0.41
1:B:233:LEU:O	1:B:237:LEU:HG	2.20	0.41
1:A:300:GLN:HB2	1:A:305:TYR:CE1	2.55	0.41
1:B:403:ARG:HG3	1:B:403:ARG:HH11	1.86	0.41
1:A:394:SER:O	1:A:398:GLN:HG3	2.21	0.41
1:A:653:ASP:HA	1:A:675:ILE:O	2.21	0.41
1:B:209:THR:HG22	1:B:209:THR:O	2.21	0.41
1:B:241:GLN:HE21	1:B:241:GLN:N	2.19	0.41
1:B:728:GLN:HE21	1:B:754:ILE:H	1.68	0.41
1:A:291:HIS:HE1	1:A:724:ASN:O	2.04	0.40
1:A:378:PHE:HA	1:A:381:SER:O	2.22	0.40
1:B:593:PHE:O	1:B:598:GLY:HA2	2.21	0.40
1:B:390:GLU:HG2	1:B:392:HIS:NE2	2.35	0.40
1:A:642:PRO:HG2	1:A:716:ILE:HG22	2.03	0.40
1:A:259:GLU:HA	1:A:260:PRO:HD2	1.89	0.40
1:B:642:PRO:HG3	1:B:743:HIS:CE1	2.57	0.40
1:B:676:THR:HG22	1:B:677:ASN:OD1	2.22	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	499/622 (80%)	458 (92%)	31 (6%)	10 (2%)	6	7
1	B	496/622 (80%)	462 (93%)	29 (6%)	5 (1%)	12	20
All	All	995/1244 (80%)	920 (92%)	60 (6%)	15 (2%)	8	12

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	513	SER
1	A	711	SER
1	B	711	SER
1	A	508	SER
1	A	512	THR
1	A	652	VAL
1	B	443	GLY
1	B	508	SER
1	A	653	ASP
1	A	209	THR
1	B	444	GLY
1	A	509	SER
1	A	712	LYS
1	A	510	PRO
1	B	636	ILE

5.3.2 Protein sidechains [i](#)

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	437/544 (80%)	416 (95%)	21 (5%)	23	40
1	B	435/544 (80%)	419 (96%)	16 (4%)	30	51
All	All	872/1088 (80%)	835 (96%)	37 (4%)	26	45

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	207	MET
1	A	213	GLU
1	A	216	ARG
1	A	227	THR
1	A	232	ARG
1	A	242	ARG
1	A	262	GLU
1	A	277	MET
1	A	385	VAL
1	A	388	SER
1	A	416	GLN
1	A	423	THR
1	A	434	LYS
1	A	436	LEU
1	A	573	GLN
1	A	622	GLN
1	A	639	GLN
1	A	641	LEU
1	A	652	VAL
1	A	703	MET
1	A	718	GLN
1	B	242	ARG
1	B	254	LEU
1	B	267	GLN
1	B	385	VAL
1	B	388	SER
1	B	416	GLN
1	B	423	THR
1	B	434	LYS
1	B	436	LEU
1	B	573	GLN
1	B	622	GLN
1	B	639	GLN
1	B	641	LEU
1	B	643	LYS

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Mol	Chain	Res	Type
1	B	703	MET
1	B	718	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	210	GLN
1	A	239	HIS
1	A	240	GLN
1	A	267	GLN
1	A	269	GLN
1	A	304	HIS
1	A	312	ASN
1	A	350	GLN
1	A	416	GLN
1	A	515	GLN
1	A	542	HIS
1	A	573	GLN
1	A	594	GLN
1	A	639	GLN
1	A	645	ASN
1	A	718	GLN
1	A	724	ASN
1	A	728	GLN
1	A	734	HIS
1	B	241	GLN
1	B	304	HIS
1	B	312	ASN
1	B	349	ASN
1	B	350	GLN
1	B	515	GLN
1	B	542	HIS
1	B	573	GLN
1	B	594	GLN
1	B	639	GLN
1	B	671	GLN
1	B	718	GLN
1	B	728	GLN
1	B	734	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.