



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:44 PM UTC

PDB ID : 3F96 / pdb_00003f96
Title : Crystal structure of human plasma platelet activating factor acetylhydrolase covalently inhibited by sarin
Authors : Samanta, U.; Bahnson, B.J.
Deposited on : 2008-11-13
Resolution : 2.10 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

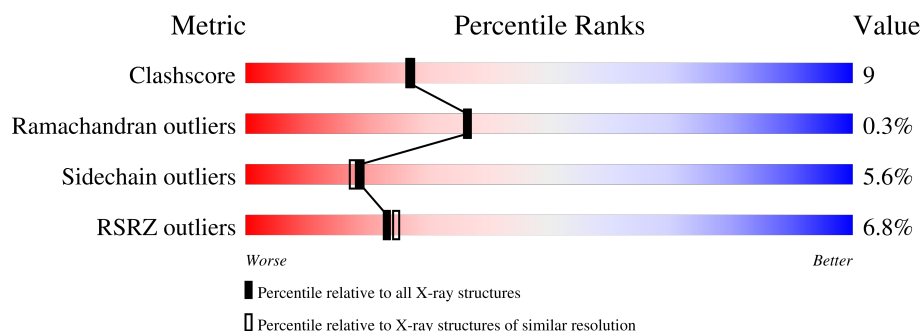
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	383	8% 68% 27% • •
1	B	383	5% 76% 19% • •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6196 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 Platelet-activating factor acetylhydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	371	Total	C	N	O	P	S	0	3	0
			3000	1919	513	554	1	13			
1	B	371	Total	C	N	O	P	S	0	3	0
			3000	1920	513	553	1	13			

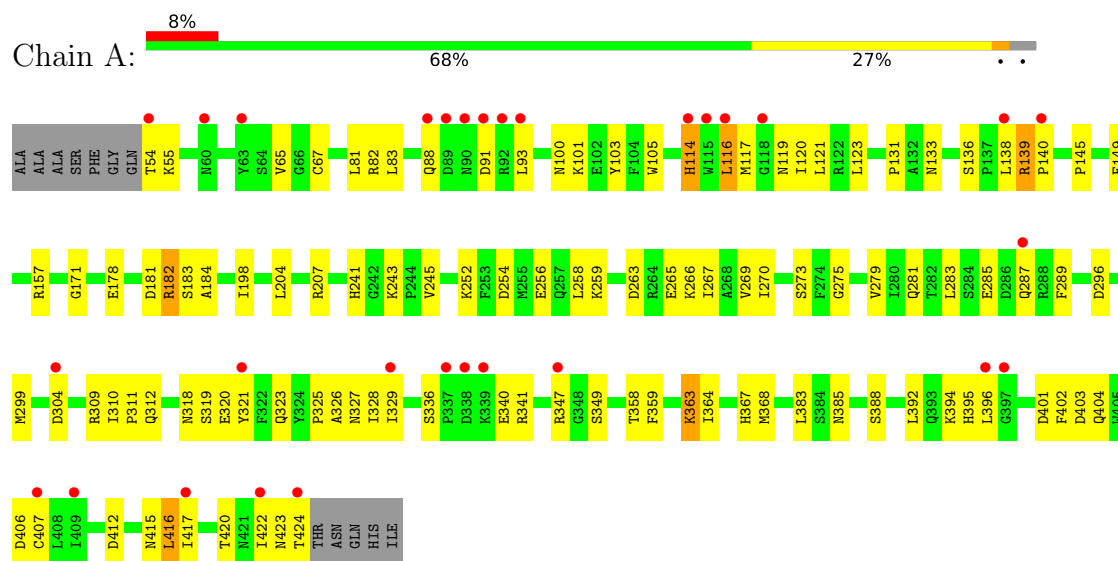
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	86	Total	O	0	0
			86	86		
2	B	110	Total	O	0	0
			110	110		

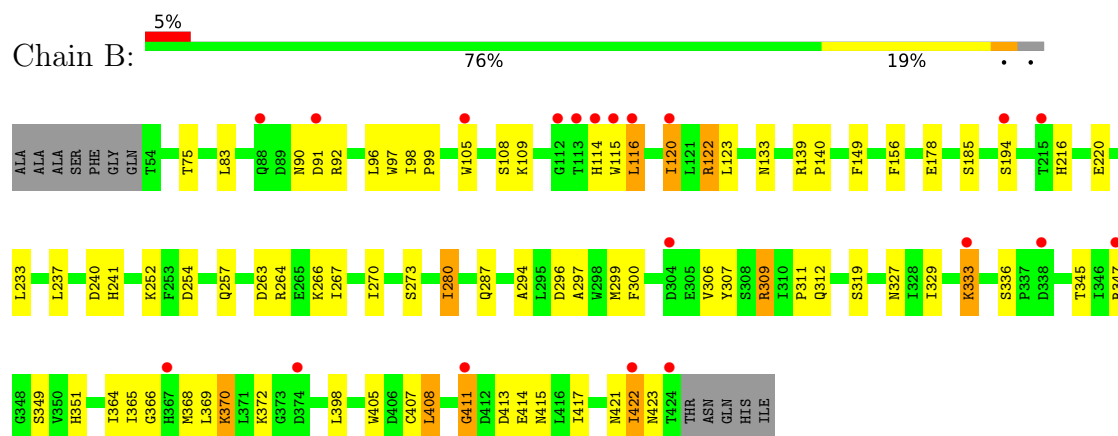
3 Residue-property plots

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: Platelet-activating factor acetylhydrolase



• Molecule 1: Platelet-activating factor acetylhydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	116.08Å 82.45Å 96.59Å 90.00° 115.11° 90.00°	Depositor
Resolution (Å)	50.00 – 2.10 50.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.7 (50.00-2.10) 95.6 (50.00-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.10Å)	Xtriage
Refinement program	REFMAC OF CCP4I FOR FINAL REFINEMENT	Depositor
R, R_{free}	0.206 , 0.278 0.210 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.321	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6196	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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: SGB

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	1.21	2/3063 (0.1%)	1.16	10/4140 (0.2%)
1	B	1.22	1/3063 (0.0%)	1.20	9/4138 (0.2%)
All	All	1.21	3/6126 (0.0%)	1.18	19/8278 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	280	ILE	CA-CB	6.54	1.62	1.54
1	A	310	ILE	CA-CB	5.78	1.61	1.54
1	A	269	VAL	CA-C	5.26	1.58	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	423	ASN	N-CA-C	10.31	126.39	112.68
1	B	411	GLY	N-CA-C	-6.97	100.86	111.24
1	B	98	ILE	N-CA-C	-6.59	101.63	108.15
1	B	413	ASP	N-CA-C	-6.18	99.33	108.79
1	A	310	ILE	CA-C-N	6.02	126.18	119.32
1	A	310	ILE	C-N-CA	6.02	126.18	119.32
1	B	300	PHE	CA-C-N	-6.00	113.44	119.56
1	B	300	PHE	C-N-CA	-6.00	113.44	119.56
1	A	323	GLN	N-CA-C	5.91	118.84	110.50
1	B	97	TRP	N-CA-C	5.78	118.07	111.02
1	A	318	ASN	N-CA-C	5.50	119.02	110.17
1	A	65	VAL	N-CA-C	5.50	115.81	108.11
1	A	321	TYR	N-CA-C	5.47	119.68	112.89
1	B	122	ARG	N-CA-CB	5.47	118.25	110.16
1	B	333	LYS	N-CA-C	-5.41	103.16	110.68
1	A	184	ALA	N-CA-C	-5.33	102.58	110.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	320	GLU	N-CA-C	5.29	116.73	111.07
1	A	402	PHE	N-CA-C	5.13	117.61	111.71
1	B	185	SER	N-CA-C	-5.03	105.69	111.07

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	3000	0	2952	67	0
1	B	3000	0	2957	47	0
2	A	86	0	0	1	0
2	B	110	0	0	1	0
All	All	6196	0	5909	110	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 9.

All (110) 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:256:GLU:HB2	1:A:259:LYS:HE2	1.49	0.93
1:B:299:MET:H	1:B:327:ASN:HD21	1.27	0.81
1:A:299:MET:H	1:A:327:ASN:HD21	1.35	0.74
1:B:422:ILE:HG22	1:B:422:ILE:O	1.87	0.74
1:A:401:ASP:O	1:A:404:GLN:HG2	1.87	0.73
1:A:182:ARG:HH11	1:A:207:ARG:CG	2.02	0.73
1:A:383:LEU:HD22	1:A:416:LEU:HD21	1.71	0.73
1:A:311:PRO:HB2	1:A:312[B]:GLN:NE2	2.05	0.71
1:A:311:PRO:HB2	1:A:312[B]:GLN:HE22	1.56	0.70
1:A:182:ARG:HH11	1:A:207:ARG:HG3	1.63	0.63
1:A:319:SER:HA	1:A:349:SER:OG	1.99	0.62
1:B:105:TRP:O	1:B:109:LYS:HG3	2.00	0.61
1:B:307:TYR:CD1	1:B:333:LYS:O	2.55	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:347:ARG:NH1	1:B:414:GLU:O	2.34	0.60
1:B:309:ARG:HH11	1:B:309:ARG:CG	2.14	0.60
1:B:307:TYR:HD1	1:B:333:LYS:O	1.86	0.59
1:A:394:LYS:HB3	1:A:395:HIS:CD2	2.39	0.58
1:A:347:ARG:HD2	1:A:417:ILE:HG12	1.86	0.58
1:A:93:LEU:HB3	1:A:131:PRO:HA	1.85	0.57
1:A:182:ARG:HH11	1:A:207:ARG:HG2	1.68	0.57
1:B:414:GLU:CD	1:B:414:GLU:H	2.12	0.57
1:B:114:HIS:CD2	1:B:115:TRP:H	2.23	0.57
1:A:392:LEU:O	1:A:396:LEU:HB2	2.04	0.56
1:B:216:HIS:NE2	1:B:220:GLU:OE2	2.38	0.56
1:A:101:LYS:HD2	1:A:105:TRP:CZ2	2.40	0.56
1:A:88:GLN:HA	1:A:139:ARG:HD2	1.87	0.56
1:A:117:MET:HE3	1:A:120:ILE:HB	1.88	0.56
1:B:311:PRO:HD2	1:B:312:GLN:OE1	2.06	0.55
1:A:265[B]:GLU:OE2	1:B:287:GLN:HG3	2.07	0.55
1:A:139:ARG:HG2	1:A:139:ARG:HH11	1.72	0.55
1:A:364:ILE:O	1:A:368:MET:HG2	2.07	0.55
1:B:90:ASN:HD22	1:B:133:ASN:HD21	1.55	0.55
1:A:309:ARG:HD3	1:B:241:HIS:CD2	2.43	0.54
1:A:182:ARG:HD3	1:A:207:ARG:HG3	1.89	0.54
1:A:326:ALA:O	1:A:329:ILE:HG22	2.08	0.54
1:B:364:ILE:O	1:B:368:MET:HG3	2.09	0.53
1:B:407:CYS:O	1:B:411:GLY:O	2.27	0.53
1:A:117:MET:HE2	1:A:121:LEU:HG	1.92	0.52
1:A:182:ARG:NH1	1:A:207:ARG:HG2	2.25	0.51
1:B:216:HIS:CE1	1:B:220:GLU:OE2	2.63	0.51
1:A:299:MET:H	1:A:327:ASN:ND2	2.06	0.51
1:A:139:ARG:HG2	1:A:139:ARG:NH1	2.25	0.51
1:A:340:GLU:C	1:A:341:ARG:HD2	2.36	0.50
1:A:114:HIS:HD2	1:A:116:LEU:H	1.58	0.50
1:A:182:ARG:HB3	1:A:207:ARG:HG3	1.94	0.50
1:A:114:HIS:CD2	1:A:116:LEU:H	2.31	0.49
1:B:91:ASP:O	1:B:92:ARG:HG3	2.12	0.49
1:A:283:LEU:HD22	1:A:312[B]:GLN:HG2	1.95	0.48
1:A:55:LYS:HB3	1:A:359:PHE:O	2.14	0.47
1:A:157:ARG:HB3	1:A:178:GLU:HB2	1.96	0.47
1:A:133:ASN:CG	1:A:136:SER:HB2	2.40	0.47
1:A:133:ASN:ND2	2:A:493:HOH:O	2.47	0.47
1:B:405:TRP:HB3	1:B:408:LEU:HD22	1.97	0.47
1:B:99:PRO:HB3	2:B:470:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:HIS:CG	1:B:115:TRP:H	2.33	0.46
1:B:254:ASP:O	1:B:257:GLN:HG3	2.15	0.46
1:A:133:ASN:HB3	1:A:136:SER:HB3	1.98	0.46
1:A:415:ASN:C	1:A:416:LEU:HD23	2.41	0.45
1:B:266:LYS:HD3	1:B:398:LEU:HD23	1.98	0.45
1:A:385:ASN:O	1:A:388:SER:HB2	2.15	0.45
1:A:287:GLN:NE2	1:B:264:ARG:HE	2.14	0.45
1:A:100:ASN:HD22	1:A:103:TYR:HE2	1.65	0.45
1:A:299:MET:N	1:A:327:ASN:HD21	2.11	0.44
1:A:275:GLY:O	1:A:279:VAL:HG23	2.18	0.43
1:A:299:MET:HE3	1:A:299:MET:HB3	1.91	0.43
1:A:263:ASP:CG	1:A:266:LYS:HD2	2.43	0.43
1:B:263:ASP:OD1	1:B:263:ASP:C	2.60	0.43
1:A:407:CYS:HB2	1:A:412:ASP:HB3	1.99	0.43
1:B:345:THR:HB	1:B:417:ILE:HB	2.00	0.43
1:B:405:TRP:O	1:B:408:LEU:HB2	2.19	0.43
1:B:90:ASN:ND2	1:B:133:ASN:HD21	2.16	0.43
1:B:139:ARG:HA	1:B:140:PRO:HD3	1.76	0.43
1:B:421:ASN:O	1:B:423:ASN:N	2.52	0.43
1:A:254:ASP:OD2	1:A:256:GLU:HG2	2.18	0.43
1:B:369:LEU:O	1:B:370:LYS:HB2	2.18	0.43
1:B:366:GLY:HA3	1:B:372:LYS:HD2	1.99	0.42
1:A:81:LEU:HD12	1:A:81:LEU:C	2.44	0.42
1:A:267:ILE:HG22	1:A:289:PHE:CD1	2.55	0.42
1:B:149:PHE:HA	1:B:270:ILE:O	2.20	0.42
1:A:139:ARG:HA	1:A:140:PRO:HD3	1.85	0.42
1:A:181:ASP:OD1	1:A:183:SER:OG	2.30	0.42
1:A:420:THR:HG23	1:A:422:ILE:O	2.20	0.42
1:B:116:LEU:O	1:B:120:ILE:HG23	2.20	0.42
1:B:319:SER:HA	1:B:349:SER:OG	2.20	0.41
1:A:138:LEU:HD22	1:A:258:LEU:HD23	2.01	0.41
1:A:241:HIS:CE1	1:B:309:ARG:HD3	2.55	0.41
1:A:363:LYS:O	1:A:367:HIS:HB2	2.19	0.41
1:B:347:ARG:HG2	1:B:417:ILE:HG13	2.01	0.41
1:A:364:ILE:O	1:A:368:MET:CG	2.68	0.41
1:B:96:LEU:HD23	1:B:99:PRO:HA	2.02	0.41
1:B:114:HIS:HD2	1:B:116:LEU:H	1.68	0.41
1:B:294:ALA:HB1	1:B:297:ALA:HB2	2.03	0.41
1:A:243:LYS:HD3	1:A:245:VAL:HG23	2.03	0.41
1:A:263:ASP:OD2	1:A:266:LYS:HD2	2.21	0.41
1:A:383:LEU:HD22	1:A:416:LEU:CD2	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:HIS:CD2	1:B:115:TRP:N	2.88	0.41
1:B:347:ARG:HG3	1:B:415:ASN:C	2.45	0.41
1:A:149:PHE:HA	1:A:270:ILE:O	2.19	0.41
1:B:237:LEU:O	1:B:240:ASP:HB3	2.21	0.41
1:A:67:CYS:SG	1:A:82:ARG:HD3	2.61	0.41
1:A:281:GLN:NE2	1:A:285:GLU:OE2	2.49	0.41
1:A:328:ILE:HG21	1:A:420:THR:HG21	2.02	0.41
1:A:403:ASP:HA	1:A:406:ASP:OD1	2.20	0.41
1:B:280:ILE:HD13	1:B:299:MET:HE1	2.01	0.41
1:B:156:PHE:HB2	1:B:178:GLU:OE2	2.22	0.40
1:B:233:LEU:HD11	1:B:267:ILE:HD13	2.02	0.40
1:A:145:PRO:HD2	1:A:171:GLY:O	2.21	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	370/383 (97%)	354 (96%)	14 (4%)	2 (0%)	24	22
1	B	370/383 (97%)	354 (96%)	16 (4%)	0	100	100
All	All	740/766 (97%)	708 (96%)	30 (4%)	2 (0%)	36	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	HIS
1	A	91	ASP

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	322/328 (98%)	304 (94%)	18 (6%)	19	18
1	B	322/328 (98%)	304 (94%)	18 (6%)	19	18
All	All	644/656 (98%)	608 (94%)	36 (6%)	19	18

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	THR
1	A	83	LEU
1	A	116	LEU
1	A	119	ASN
1	A	123	LEU
1	A	139	ARG
1	A	182	ARG
1	A	198	ILE
1	A	204	LEU
1	A	252	LYS
1	A	296	ASP
1	A	304	ASP
1	A	325	PRO
1	A	336	SER
1	A	358	THR
1	A	363	LYS
1	A	416	LEU
1	A	424	THR
1	B	75	THR
1	B	83	LEU
1	B	108	SER
1	B	116	LEU
1	B	120	ILE
1	B	122	ARG
1	B	123	LEU
1	B	194	SER
1	B	252	LYS

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Mol	Chain	Res	Type
1	B	296	ASP
1	B	306	VAL
1	B	309	ARG
1	B	329	ILE
1	B	336	SER
1	B	365	ILE
1	B	370	LYS
1	B	408	LEU
1	B	422	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	100	ASN
1	A	114	HIS
1	A	133	ASN
1	A	135	ASN
1	A	327	ASN
1	A	393	GLN
1	A	404	GLN
1	A	415	ASN
1	B	90	ASN
1	B	114	HIS
1	B	135	ASN
1	B	231	GLN
1	B	287	GLN
1	B	327	ASN
1	B	423	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
1	SGB	B	273[A]	-	10,12,13	3.97	3 (30%)	8,16,18	2.71	3 (37%)
1	SGB	B	273[B]	-	10,12,13	3.93	3 (30%)	8,16,18	2.59	3 (37%)
1	SGB	A	273[A]	-	10,12,13	4.30	3 (30%)	8,16,18	1.66	2 (25%)
1	SGB	A	273[B]	-	10,12,13	4.27	3 (30%)	8,16,18	1.67	1 (12%)

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
1	SGB	B	273[A]	-	-	1/11/13/15	-
1	SGB	B	273[B]	-	-	3/11/13/15	-
1	SGB	A	273[A]	-	-	1/11/13/15	-
1	SGB	A	273[B]	-	-	4/11/13/15	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	273[A]	SGB	P1-O1	9.35	1.61	1.47
1	B	273[B]	SGB	P1-O1	9.35	1.61	1.47
1	A	273[A]	SGB	P1-O1	8.86	1.61	1.47
1	A	273[B]	SGB	P1-O1	8.86	1.61	1.47
1	A	273[A]	SGB	P1-OG	8.53	1.68	1.58
1	A	273[B]	SGB	P1-OG	8.53	1.68	1.58
1	B	273[A]	SGB	P1-OG	6.15	1.65	1.58
1	B	273[B]	SGB	P1-OG	6.15	1.65	1.58
1	B	273[A]	SGB	O-C	5.08	1.39	1.20
1	B	273[B]	SGB	O-C	5.08	1.39	1.20
1	A	273[A]	SGB	O-C	4.92	1.38	1.20
1	A	273[B]	SGB	O-C	4.92	1.38	1.20

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	273[A]	SGB	OG-CB-CA	5.91	113.90	108.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	273[B]	SGB	OG-CB-CA	5.91	113.90	108.14
1	B	273[A]	SGB	O1-P1-C1	-3.48	102.63	114.97
1	B	273[B]	SGB	O1-P1-C1	-3.25	103.44	114.97
1	A	273[B]	SGB	O1-P1-C1	-3.12	103.90	114.97
1	A	273[A]	SGB	OG-P1-C1	2.69	112.40	105.02
1	A	273[A]	SGB	O1-P1-C1	-2.57	105.85	114.97
1	B	273[A]	SGB	OG-P1-O1	2.09	118.68	113.15
1	B	273[B]	SGB	OG-P1-O1	2.09	118.68	113.15

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	273[B]	SGB	C2-O2-P1-O1
1	A	273[B]	SGB	C2-O2-P1-O1
1	B	273[B]	SGB	CB-OG-P1-O2
1	A	273[B]	SGB	C4-C2-O2-P1
1	A	273[A]	SGB	N-CA-CB-OG
1	A	273[B]	SGB	N-CA-CB-OG
1	B	273[A]	SGB	N-CA-CB-OG
1	B	273[B]	SGB	N-CA-CB-OG
1	A	273[B]	SGB	C3-C2-O2-P1

There are no ring outliers.

No monomer is involved in short contacts.

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

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	370/383 (96%)	0.71	30 (8%) 18 19	18, 38, 58, 84	2 (0%)
1	B	370/383 (96%)	0.20	20 (5%) 31 33	17, 31, 52, 74	2 (0%)
All	All	740/766 (96%)	0.46	50 (6%) 23 25	17, 34, 56, 84	4 (0%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	89	ASP	6.3
1	A	54	THR	5.5
1	B	115	TRP	4.7
1	A	417	ILE	4.6
1	A	115	TRP	4.4
1	A	92	ARG	4.2
1	A	424	THR	4.0
1	B	411	GLY	3.9
1	A	93	LEU	3.8
1	A	329	ILE	3.6
1	A	339	LYS	3.5
1	A	396	LEU	3.3
1	B	113	THR	3.0
1	B	114	HIS	3.0
1	A	304	ASP	2.8
1	A	114	HIS	2.8
1	A	90	ASN	2.7
1	A	91	ASP	2.7
1	A	397	GLY	2.6
1	A	116	LEU	2.6
1	A	88	GLN	2.6
1	A	338	ASP	2.5
1	B	91	ASP	2.5
1	B	88	GLN	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	118	GLY	2.5
1	A	409	ILE	2.5
1	B	120	ILE	2.5
1	A	287	GLN	2.4
1	B	194	SER	2.4
1	A	337	PRO	2.4
1	B	367	HIS	2.4
1	A	407	CYS	2.4
1	B	422	ILE	2.3
1	B	338	ASP	2.3
1	B	116	LEU	2.3
1	A	138	LEU	2.3
1	B	112	GLY	2.2
1	B	105	TRP	2.2
1	A	321	TYR	2.2
1	B	347	ARG	2.2
1	A	63	TYR	2.2
1	A	140	PRO	2.1
1	B	333	LYS	2.1
1	A	60	ASN	2.1
1	B	304	ASP	2.1
1	A	422	ILE	2.1
1	B	374	ASP	2.0
1	A	347	ARG	2.0
1	B	424	THR	2.0
1	B	215	THR	2.0

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

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
1	SGB	B	273[A]	13/14	0.67	0.24	21,37,50,54	5
1	SGB	B	273[B]	13/14	0.67	0.24	21,37,52,54	5
1	SGB	A	273[A]	13/14	0.68	0.26	26,38,52,56	5
1	SGB	A	273[B]	13/14	0.68	0.26	26,38,53,56	5

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.