



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 12:29 AM UTC

PDB ID : 1EJD / pdb_00001ejd
Title : Crystal structure of unliganded mura (type1)
Authors : Eschenburg, S.; Schonbrunn, E.
Deposited on : 2000-03-02
Resolution : 1.55 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

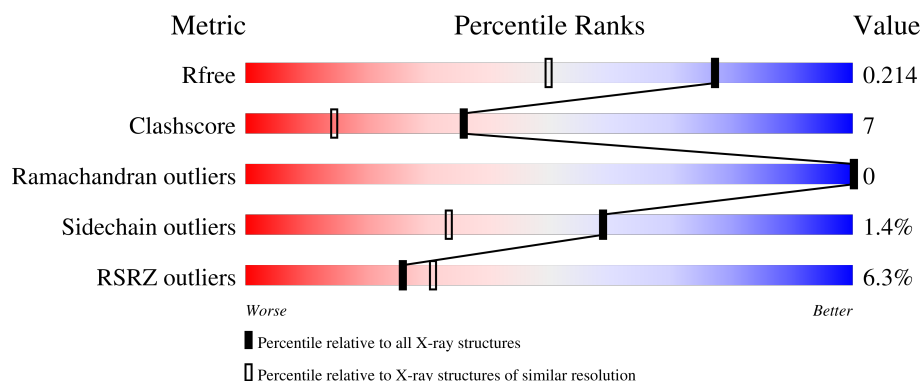
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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	419	4% 83% 16%
1	B	419	9% 84% 15%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7522 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.

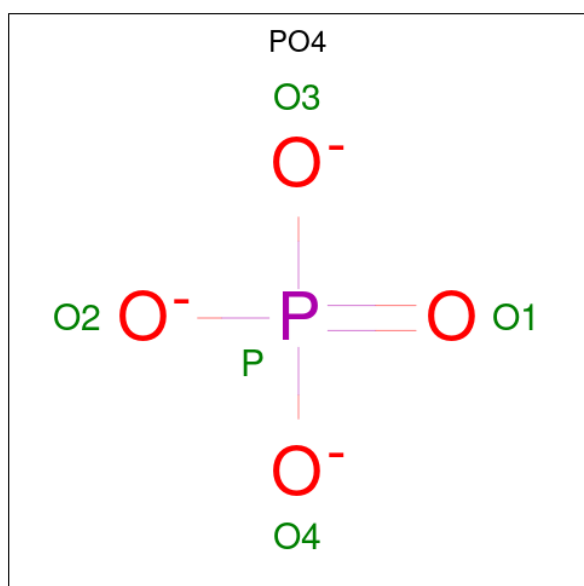
- Molecule 1 is a protein called UDP-N-ACETYLGLUCOSAMINE ENOLPYRUVYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	419	Total	C	N	O	S	1	21	0
			3239	2044	573	606	16			
1	B	419	Total	C	N	O	S	0	24	0
			3250	2051	577	606	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	67	IAS	ASN	conflict	UNP P33038
B	67	IAS	ASN	conflict	UNP P33038

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



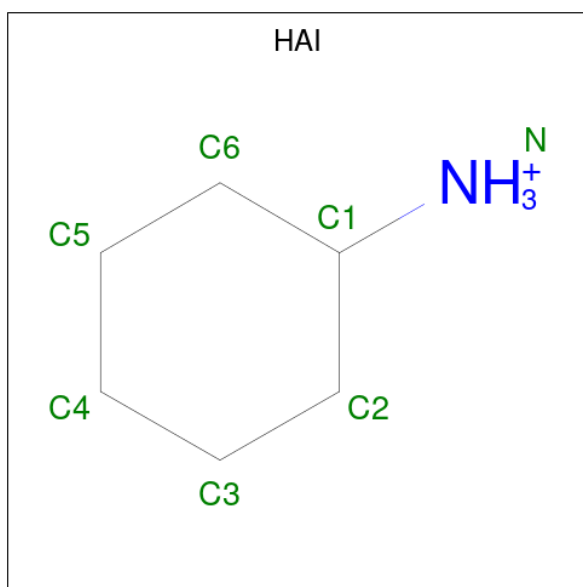
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		

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

- Molecule 3 is CYCLOHEXYLAMMONIUM ION (CCD ID: HAI) (formula: C₆H₁₄N).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			7	6	1		
3	A	1	Total	C	N	0	0
			7	6	1		
3	B	1	Total	C	N	0	0
			7	6	1		
3	B	1	Total	C	N	0	0
			7	6	1		
3	B	1	Total	C	N	0	0
			7	6	1		

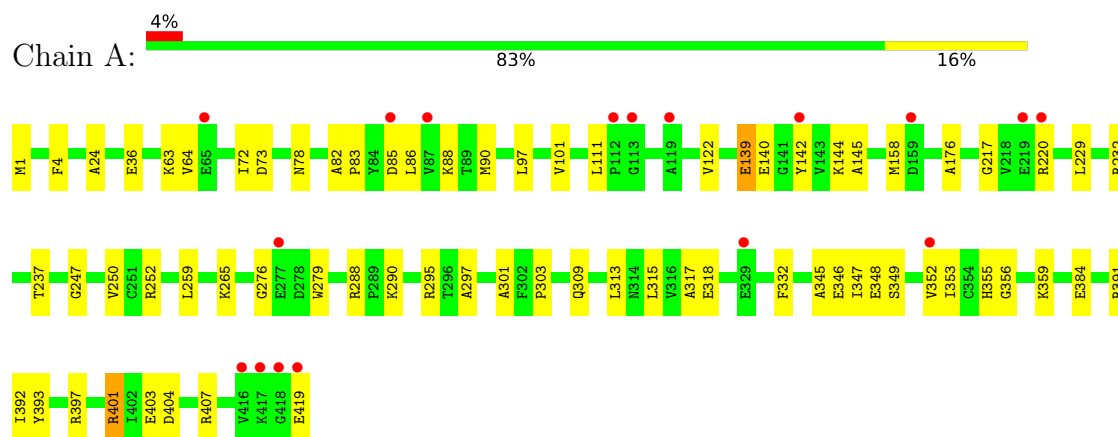
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	480	Total	O	0	2
			480	480		
4	B	443	Total	O	0	3
			443	443		

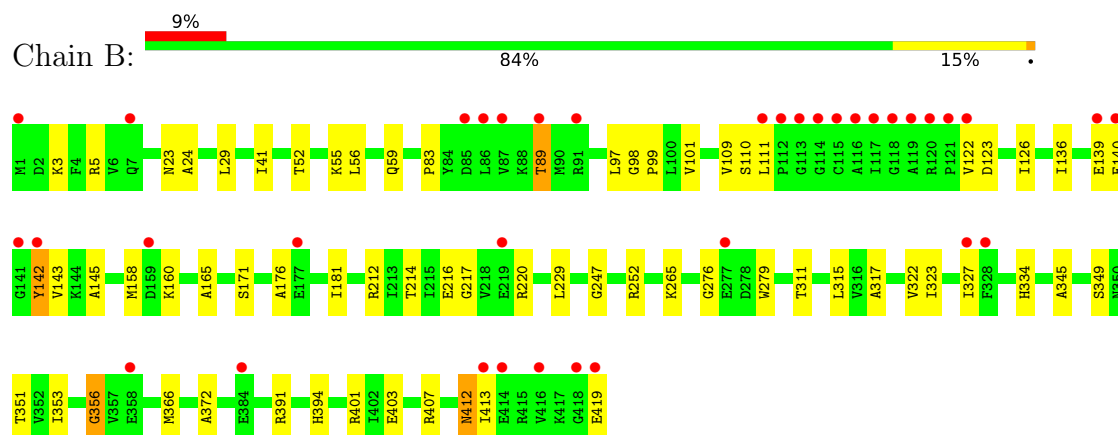
3 Residue-property plots

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: UDP-N-ACETYLGLUCOSAMINE ENOLPYRUVYLTRANSFERASE



• Molecule 1: UDP-N-ACETYLGLUCOSAMINE ENOLPYRUVYLTRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	87.18Å 156.30Å 84.04Å 90.00° 91.63° 90.00°	Depositor
Resolution (Å)	17.00 – 1.55 17.00 – 1.55	Depositor EDS
% Data completeness (in resolution range)	96.0 (17.00-1.55) 95.8 (17.00-1.55)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 1.55Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.185 , 0.207 0.194 , 0.214	Depositor DCC
R_{free} test set	4686 reflections (3.01%)	wwPDB-VP
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 45.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.007 for -1/2*h-1/2*k,-3/2*h+1/2*k,-l 0.007 for -1/2*h+1/2*k,3/2*h+1/2*k,-l 0.002 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.004 for 1/2*h+1/2*k,3/2*h-1/2*k,-l 0.015 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7522	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% 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: IAS, HAI, 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	0.55	0/3359	0.96	11/4543 (0.2%)
1	B	0.53	0/3382	0.93	7/4574 (0.2%)
All	All	0.54	0/6741	0.94	18/9117 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	24	ALA	N-CA-C	-7.96	103.53	113.55
1	A	82	ALA	CA-C-N	7.80	127.72	119.85
1	A	82	ALA	C-N-CA	7.80	127.72	119.85
1	B	356	GLY	N-CA-C	6.36	118.80	111.36
1	A	247	GLY	N-CA-C	6.06	121.95	112.31
1	B	97	LEU	N-CA-C	6.02	117.52	111.07
1	A	349	SER	CB-CA-C	-5.78	109.93	116.63
1	B	349	SER	CB-CA-C	-5.59	110.15	116.63
1	A	24	ALA	N-CA-C	-5.54	106.56	113.55
1	A	90	MET	N-CA-C	5.41	117.02	107.61
1	A	139	GLU	N-CA-C	-5.32	100.99	109.23
1	A	288	ARG	CA-C-N	-5.31	115.11	120.31
1	A	288	ARG	C-N-CA	-5.31	115.11	120.31
1	B	123	ASP	N-CA-C	5.30	117.47	111.11
1	A	303	PRO	N-CA-C	5.14	118.93	110.55
1	B	142	TYR	N-CA-C	5.13	115.93	108.14
1	A	88	LYS	N-CA-C	5.11	117.59	111.71
1	B	247	GLY	N-CA-C	5.09	120.05	112.41

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	3239	0	3337	39	0
1	B	3250	0	3354	59	1
2	A	35	0	0	1	0
2	B	40	0	0	0	0
3	A	14	0	28	1	0
3	B	21	0	42	2	0
4	A	480	0	0	3	1
4	B	443	0	0	14	2
All	All	7522	0	6761	98	4

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 7.

All (98) 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:252[A]:ARG:CZ	4:B:427[A]:HOH:O	1.92	1.15
1:B:252[A]:ARG:NH2	4:B:427[A]:HOH:O	2.01	0.86
1:B:101[B]:VAL:HG21	1:B:145:ALA:HB3	1.63	0.79
1:A:347:ILE:HD13	1:A:352:VAL:HG22	1.66	0.77
1:A:250:VAL:HG11	3:B:1420:HAI:H32	1.71	0.72
1:B:126:ILE:HD12	1:B:143:VAL:HG21	1.72	0.71
1:B:216:GLU:HG3	4:B:458:HOH:O	1.91	0.70
1:B:276:GLY:HA3	1:B:279:TRP:CE2	2.29	0.67
1:A:346:GLU:OE1	1:A:355:HIS:HE1	1.78	0.67
1:B:412:ASN:C	1:B:412:ASN:HD22	2.03	0.67
1:B:212[B]:ARG:NH2	4:B:596:HOH:O	2.29	0.65
1:B:322[A]:VAL:HG13	1:B:351[A]:THR:CG2	2.27	0.65
1:B:394:HIS:HB2	4:B:632:HOH:O	1.98	0.63
1:B:98:GLY:HA3	4:B:841:HOH:O	2.00	0.62
1:B:322[A]:VAL:HG13	1:B:351[A]:THR:HG23	1.83	0.61
1:A:252[A]:ARG:HH21	2:A:2433:PO4:P	2.26	0.59
1:A:359:LYS:HD2	1:A:384:GLU:CD	2.28	0.59
1:B:29:LEU:HD23	1:B:41[B]:ILE:HD12	1.85	0.58
1:A:401:ARG:HG2	1:A:401:ARG:HH11	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101[B]:VAL:HG21	1:B:145:ALA:CB	2.31	0.58
1:A:36:GLU:OE1	1:A:220:ARG:HD3	2.04	0.58
1:B:158[A]:MET:HE1	1:B:165:ALA:HB3	1.86	0.56
1:B:181[B]:ILE:CD1	1:B:214:THR:HG23	2.35	0.56
1:B:3:LYS:HE3	1:B:419:GLU:HA	1.86	0.56
1:A:237:THR:HA	1:A:309[B]:GLN:OE1	2.06	0.55
1:A:85:ASP:OD1	1:A:86:LEU:HG	2.07	0.55
1:B:322[A]:VAL:HG22	1:B:353:ILE:HA	1.88	0.55
1:B:126:ILE:CG2	1:B:136:ILE:HG21	2.37	0.55
1:A:139:GLU:HB2	1:A:142:TYR:CE2	2.42	0.54
1:A:297:ALA:HB3	1:A:301:ALA:CB	2.37	0.54
1:B:322[A]:VAL:CG1	1:B:351[A]:THR:HG23	2.37	0.54
1:B:126:ILE:HG23	1:B:136:ILE:HG21	1.90	0.54
1:A:397[A]:ARG:HH21	1:A:397[A]:ARG:HG3	1.72	0.54
1:B:99:PRO:HD3	4:B:841:HOH:O	2.07	0.54
1:B:5[B]:ARG:HD3	4:B:625:HOH:O	2.07	0.54
1:B:311:THR:HA	1:B:323:ILE:HD12	1.90	0.53
1:A:122[A]:VAL:HG22	4:A:851:HOH:O	2.09	0.53
1:B:276:GLY:HA3	1:B:279:TRP:NE1	2.24	0.53
1:B:403:GLU:O	1:B:407:ARG:HG3	2.09	0.53
1:B:29:LEU:HD23	1:B:41[B]:ILE:CD1	2.39	0.52
1:A:111:LEU:HD22	1:A:122[A]:VAL:HG11	1.90	0.52
1:B:322[A]:VAL:CG1	1:B:351[A]:THR:CG2	2.87	0.52
1:A:101[A]:VAL:HG11	1:A:145:ALA:HB3	1.92	0.52
1:B:52:THR:HA	1:B:55:LYS:HE3	1.92	0.52
1:B:334:HIS:HB3	1:B:372:ALA:HB1	1.93	0.50
1:B:401:ARG:HB2	4:B:731:HOH:O	2.11	0.50
1:B:171[B]:SER:OG	4:B:841:HOH:O	2.17	0.50
1:B:126:ILE:CD1	1:B:143:VAL:HG21	2.40	0.50
1:A:348:GLU:CD	1:A:353:ILE:HD11	2.37	0.49
1:B:252[A]:ARG:NH1	4:B:427[A]:HOH:O	2.23	0.49
1:B:59[B]:GLN:HE22	1:B:83:PRO:HD3	1.78	0.49
1:A:97:LEU:O	1:A:101[B]:VAL:HG23	2.13	0.49
1:B:322[A]:VAL:HG22	1:B:353:ILE:HG12	1.95	0.48
1:B:279:TRP:CD1	3:B:1420:HAI:H31	2.49	0.48
1:A:295[A]:ARG:HD2	4:A:526:HOH:O	2.14	0.47
1:A:317:ALA:O	1:A:356:GLY:HA3	2.15	0.47
1:B:142:TYR:CD1	1:B:142:TYR:C	2.93	0.47
1:B:23:ASN:ND2	4:B:778:HOH:O	2.47	0.46
1:B:56:LEU:C	1:B:56:LEU:HD23	2.40	0.46
1:B:412:ASN:C	1:B:412:ASN:ND2	2.69	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:VAL:HG13	1:A:72[A]:ILE:HD13	1.98	0.46
1:B:176:ALA:O	1:B:217:GLY:HA3	2.16	0.45
1:B:101[B]:VAL:CG2	1:B:145:ALA:HB3	2.42	0.45
1:A:232:ARG:O	1:A:259[A]:LEU:HD21	2.17	0.45
1:A:315:LEU:HD21	1:A:345:ALA:HB2	1.99	0.45
1:B:317:ALA:O	1:B:356:GLY:HA3	2.16	0.45
1:A:276:GLY:HA3	1:A:279:TRP:CE2	2.51	0.44
1:B:109[B]:VAL:HG12	1:B:110:SER:O	2.18	0.44
1:B:111:LEU:HD22	1:B:122[B]:VAL:HG21	2.00	0.44
1:B:252[B]:ARG:HB2	1:B:252[B]:ARG:NH1	2.31	0.44
1:A:290:LYS:NZ	1:A:318:GLU:OE1	2.42	0.44
1:B:139:GLU:O	1:B:140:GLU:C	2.60	0.43
1:A:176:ALA:O	1:A:217:GLY:HA3	2.18	0.43
1:B:315:LEU:HD21	1:B:345:ALA:HB2	2.00	0.43
1:A:4:PHE:CD2	1:A:392:ILE:HG21	2.53	0.43
1:B:366[B]:MET:SD	1:B:391:ARG:HD3	2.58	0.43
1:B:5[A]:ARG:HD2	4:B:756:HOH:O	2.18	0.43
1:A:403:GLU:HG2	1:A:404:ASP:N	2.33	0.43
1:A:407:ARG:HD2	4:A:839:HOH:O	2.18	0.42
1:B:111:LEU:HD22	1:B:122[A]:VAL:HG11	2.00	0.42
1:A:313:LEU:C	1:A:313:LEU:HD23	2.44	0.42
1:A:297:ALA:HB3	1:A:301:ALA:HB3	2.02	0.41
1:A:63:LYS:HB2	1:A:73:ASP:HB3	2.02	0.41
1:B:160:LYS:HD3	1:B:160:LYS:HA	1.74	0.41
1:A:1:MET:HE3	1:A:391:ARG:NH1	2.36	0.41
1:B:327:ILE:HD11	4:B:874:HOH:O	2.20	0.41
1:A:158[B]:MET:HE2	1:A:158[B]:MET:HB3	1.98	0.41
1:B:89:THR:O	1:B:89:THR:OG1	2.38	0.41
1:B:158[A]:MET:HE2	1:B:158[A]:MET:HB3	1.96	0.41
1:A:144:LYS:HE3	3:A:1424:HAI:H62	2.03	0.40
1:A:97:LEU:O	1:A:101[A]:VAL:HG22	2.21	0.40
1:A:393:TYR:O	1:A:397[B]:ARG:HG3	2.22	0.40
1:B:252[B]:ARG:CZ	1:B:252[B]:ARG:CB	2.99	0.40
1:B:252[B]:ARG:CZ	1:B:252[B]:ARG:HB3	2.51	0.40
1:A:139:GLU:O	1:A:140:GLU:C	2.63	0.40
1:A:332:PHE:CE1	1:A:352:VAL:HG23	2.56	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:757:HOH:O	4:B:757:HOH:O[2_655]	1.59	0.61
4:B:797:HOH:O	4:B:797:HOH:O[2_656]	1.66	0.54
1:B:220:ARG:NH2	1:B:220:ARG:NH2[2_656]	1.99	0.21
4:A:426:HOH:O	4:A:426:HOH:O[2_556]	2.08	0.12

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	436/419 (104%)	429 (98%)	7 (2%)	0	100	100
1	B	439/419 (105%)	437 (100%)	2 (0%)	0	100	100
All	All	875/838 (104%)	866 (99%)	9 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	350/329 (106%)	345 (99%)	5 (1%)	59	33
1	B	353/329 (107%)	349 (99%)	4 (1%)	65	43
All	All	703/658 (107%)	694 (99%)	9 (1%)	59	36

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	ASN
1	A	83	PRO
1	A	229	LEU
1	A	401	ARG
1	A	419	GLU
1	B	89	THR
1	B	229	LEU
1	B	412	ASN
1	B	413	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	59	GLN
1	A	78	ASN
1	A	148	ASN
1	A	330	ASN
1	A	355	HIS
1	B	76	ASN
1	B	108	GLN
1	B	148	ASN
1	B	412	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	IAS	B	67	1	6,7,8	1.03	0	3,8,10	1.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	IAS	A	67	1	6,7,8	0.92	0	3,8,10	1.13	0

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	IAS	B	67	1	-	0/7/7/8	-
1	IAS	A	67	1	-	0/7/7/8	-

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.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

20 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	HAI	B	1421	-	7,7,7	1.28	1 (14%)	8,8,8	1.59	2 (25%)
2	PO4	B	2439	-	4,4,4	1.67	1 (25%)	6,6,6	0.51	0
2	PO4	B	2427[A]	-	4,4,4	1.67	0	6,6,6	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	A	2430	-	4,4,4	1.68	0	6,6,6	0.46	0
2	PO4	A	2436	-	4,4,4	1.65	0	6,6,6	0.46	0
2	PO4	A	2432[A]	-	4,4,4	1.71	0	6,6,6	0.45	0
3	HAI	B	1422	-	7,7,7	1.03	0	8,8,8	1.76	2 (25%)
2	PO4	B	2429	-	4,4,4	1.77	2 (50%)	6,6,6	0.46	0
2	PO4	B	2426	-	4,4,4	1.68	0	6,6,6	0.47	0
2	PO4	A	2437	-	4,4,4	1.65	1 (25%)	6,6,6	0.45	0
2	PO4	B	2425	-	4,4,4	1.74	1 (25%)	6,6,6	0.47	0
2	PO4	A	2433	-	4,4,4	1.78	2 (50%)	6,6,6	0.49	0
2	PO4	B	2424[A]	-	4,4,4	1.72	1 (25%)	6,6,6	0.45	0
3	HAI	B	1420	-	7,7,7	1.32	1 (14%)	8,8,8	2.33	2 (25%)
3	HAI	A	1423	-	7,7,7	1.32	1 (14%)	8,8,8	1.41	2 (25%)
2	PO4	A	2431	-	4,4,4	1.45	0	6,6,6	0.42	0
3	HAI	A	1424	-	7,7,7	1.15	0	8,8,8	2.00	3 (37%)
2	PO4	A	2434	-	4,4,4	1.76	1 (25%)	6,6,6	0.48	0
2	PO4	B	2428	-	4,4,4	1.77	1 (25%)	6,6,6	0.47	0
2	PO4	B	2422	-	4,4,4	1.52	0	6,6,6	0.48	0

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	HAI	B	1420	-	-	-	0/1/1/1
3	HAI	A	1423	-	-	-	0/1/1/1
3	HAI	B	1422	-	-	-	0/1/1/1
3	HAI	B	1421	-	-	-	0/1/1/1
3	HAI	A	1424	-	-	-	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1420	HAI	C5-C6	2.95	1.60	1.53
3	A	1423	HAI	C5-C6	2.48	1.59	1.53
3	B	1421	HAI	C5-C6	2.47	1.59	1.53
2	B	2429	PO4	P-O3	-2.08	1.48	1.54
2	A	2433	PO4	P-O2	-2.08	1.48	1.54
2	B	2424[A]	PO4	P-O2	-2.07	1.48	1.54
2	B	2439	PO4	P-O4	-2.07	1.48	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2428	PO4	P-O4	-2.07	1.48	1.54
2	B	2429	PO4	P-O2	-2.04	1.48	1.54
2	A	2437	PO4	P-O2	-2.04	1.48	1.54
2	A	2434	PO4	P-O2	-2.04	1.48	1.54
2	B	2425	PO4	P-O3	-2.02	1.48	1.54
2	A	2433	PO4	P-O3	-2.02	1.48	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1420	HAI	C6-C1-C2	4.93	115.30	110.29
3	A	1424	HAI	C6-C1-C2	3.57	113.92	110.29
3	B	1420	HAI	C5-C6-C1	3.33	116.28	111.77
3	B	1422	HAI	C6-C1-C2	3.26	113.60	110.29
3	A	1424	HAI	C6-C1-N	3.04	120.21	111.17
3	B	1421	HAI	C6-C1-C2	2.74	113.07	110.29
3	A	1423	HAI	C6-C1-C2	2.55	112.88	110.29
3	A	1424	HAI	C2-C1-N	2.54	118.72	111.17
3	B	1421	HAI	C5-C6-C1	2.44	115.07	111.77
3	B	1422	HAI	C2-C1-N	2.40	118.33	111.17
3	A	1423	HAI	C5-C6-C1	2.39	115.01	111.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2433	PO4	1	0
3	B	1420	HAI	2	0
3	A	1424	HAI	1	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	418/419 (99%)	0.28	17 (4%) 41 49	12, 23, 36, 62	22 (5%)
1	B	418/419 (99%)	0.37	36 (8%) 16 18	11, 23, 41, 64	24 (5%)
All	All	836/838 (99%)	0.32	53 (6%) 26 31	11, 23, 38, 64	46 (5%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	113	GLY	7.0
1	A	418	GLY	5.9
1	B	115	CYS	5.5
1	B	419	GLU	5.3
1	B	119	ALA	5.2
1	B	116	ALA	5.1
1	A	142	TYR	4.7
1	B	142	TYR	4.6
1	B	121	PRO	4.2
1	A	419	GLU	4.0
1	B	141	GLY	3.9
1	A	85	ASP	3.7
1	B	89	THR	3.5
1	B	113	GLY	3.5
1	B	416	VAL	3.5
1	B	111	LEU	3.1
1	B	1	MET	3.1
1	A	65	GLU	3.0
1	B	418	GLY	2.9
1	B	159	ASP	2.9
1	B	413	ILE	2.8
1	A	329	GLU	2.8
1	B	86	LEU	2.8
1	B	112	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	118	GLY	2.7
1	B	85	ASP	2.7
1	B	140	GLU	2.7
1	B	87	VAL	2.6
1	B	114	GLY	2.6
1	B	384	GLU	2.5
1	A	159	ASP	2.5
1	A	112	PRO	2.5
1	B	328	PHE	2.4
1	B	122[A]	VAL	2.4
1	A	119	ALA	2.4
1	B	7	GLN	2.4
1	B	139	GLU	2.4
1	B	120	ARG	2.4
1	B	277	GLU	2.4
1	B	91	ARG	2.3
1	A	416	VAL	2.3
1	A	277	GLU	2.3
1	B	117	ILE	2.2
1	A	219	GLU	2.2
1	A	417	LYS	2.2
1	B	327	ILE	2.1
1	B	219	GLU	2.1
1	A	87	VAL	2.1
1	A	352	VAL	2.1
1	A	220	ARG	2.1
1	B	358	GLU	2.1
1	B	414	GLU	2.1
1	B	177	GLU	2.0

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

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
1	IAS	B	67	8/9	0.80	0.14	29,31,35,35	0
1	IAS	A	67	8/9	0.86	0.14	28,29,34,34	0

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	PO4	B	2427[A]	5/5	0.63	0.19	47,47,48,49	5
3	HAI	A	1423	7/7	0.67	0.17	58,58,58,59	0
2	PO4	B	2439	5/5	0.69	0.21	33,34,38,40	5
2	PO4	B	2426	5/5	0.70	0.19	52,52,53,54	5
2	PO4	A	2433	5/5	0.72	0.20	37,38,42,44	5
2	PO4	A	2432[A]	5/5	0.74	0.19	52,52,53,53	5
2	PO4	A	2437	5/5	0.76	0.15	31,36,39,39	5
2	PO4	B	2422	5/5	0.76	0.15	32,36,36,37	5
2	PO4	B	2424[A]	5/5	0.78	0.14	47,49,51,52	5
2	PO4	A	2436	5/5	0.79	0.14	41,41,42,43	5
3	HAI	B	1421	7/7	0.79	0.14	55,56,56,56	0
3	HAI	B	1420	7/7	0.83	0.17	44,45,46,47	0
2	PO4	B	2425	5/5	0.84	0.15	28,28,31,31	5
3	HAI	B	1422	7/7	0.85	0.17	46,47,48,49	0
2	PO4	B	2428	5/5	0.87	0.16	41,41,42,42	5
2	PO4	A	2434	5/5	0.88	0.13	44,45,46,47	5
2	PO4	A	2430	5/5	0.88	0.15	55,56,56,58	5
2	PO4	B	2429	5/5	0.89	0.14	40,40,41,41	5
2	PO4	A	2431	5/5	0.89	0.15	25,28,33,34	5
3	HAI	A	1424	7/7	0.90	0.11	33,35,36,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.