



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:48 AM UTC

PDB ID : 2HHF / pdb_00002hhf
Title : X-ray crystal structure of oxidized human mitochondrial branched chain aminotransferase (hBCATm)
Authors : Yennawar, N.H.; Hutson, S.M.
Deposited on : 2006-06-28
Resolution : 1.80 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	:	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

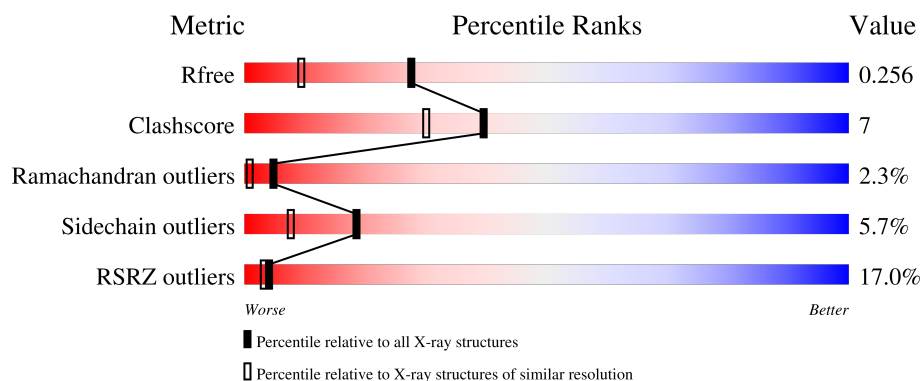
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	365	16% 75% 16% 5%
2	B	365	17% 84% 12% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5994 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 Branched-chain-amino-acid aminotransferase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	348	2799	1795	488	498	18	0	3	1

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	108	OCS	CYS	modified residue	UNP O15382
A	141	TYO	TYR	modified residue	UNP O15382
A	159	ARG	THR	conflict	UNP O15382

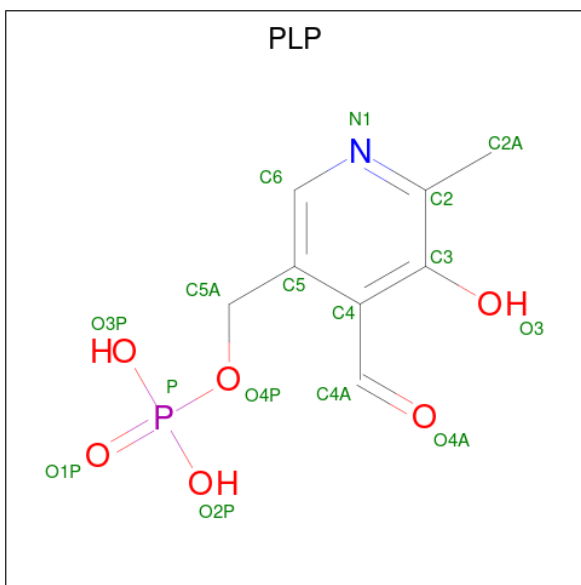
- Molecule 2 is a protein called Branched-chain-amino-acid aminotransferase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	365	2944	1893	513	520	18	0	4	0

There are 2 discrepancies between the modelled and reference sequences:

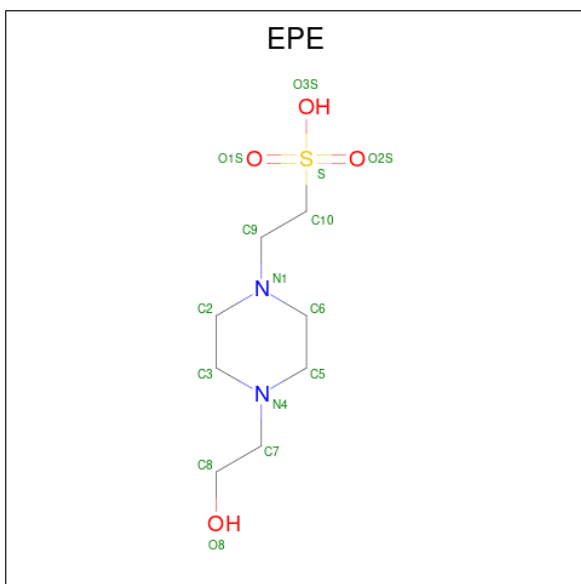
Chain	Residue	Modelled	Actual	Comment	Reference
B	608	OCS	CYS	modified residue	UNP O15382
B	659	ARG	THR	conflict	UNP O15382

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: C₈H₁₀NO₆P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

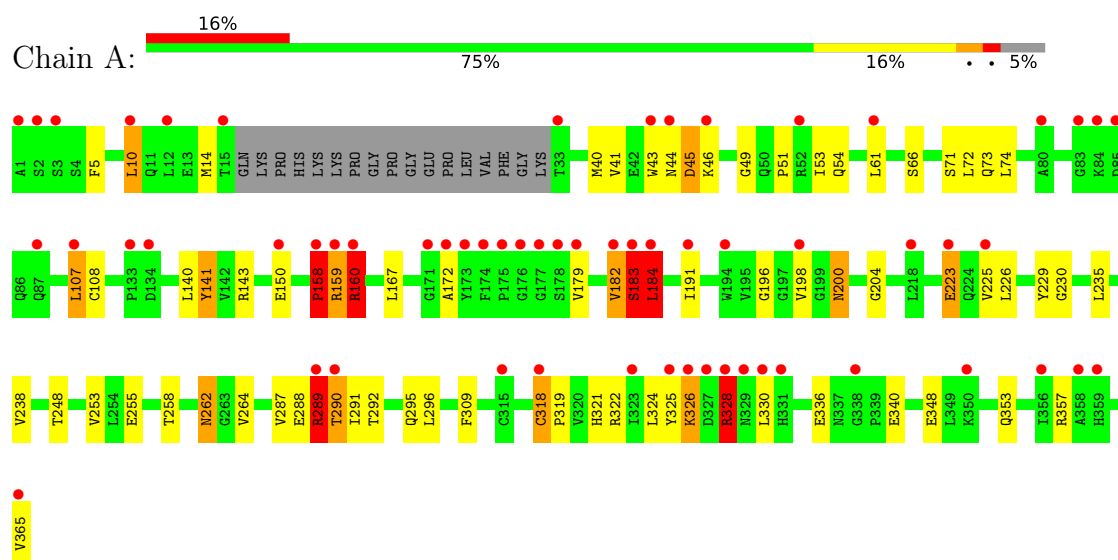
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	97	Total 97	O 97	0	0
5	B	109	Total 109	O 109	0	0

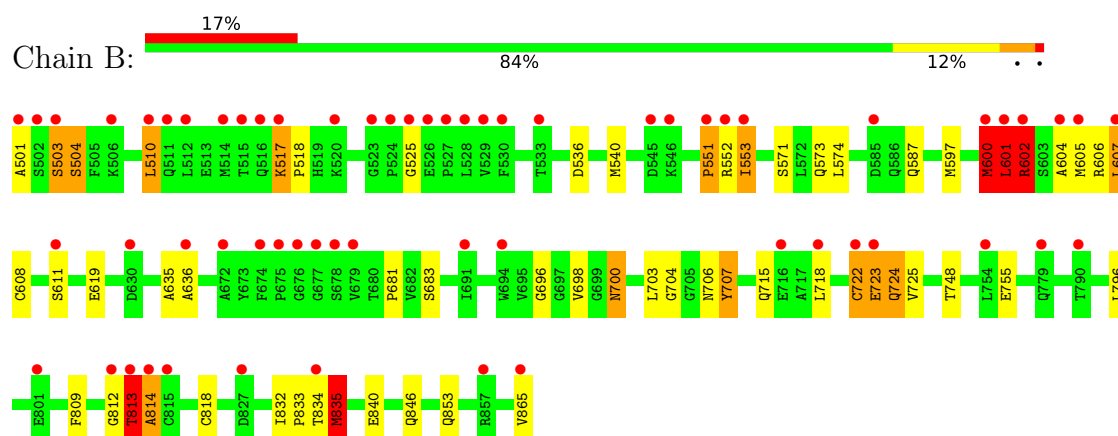
3 Residue-property plots

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: Branched-chain-amino-acid aminotransferase, mitochondrial



- Molecule 2: Branched-chain-amino-acid aminotransferase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.12Å 104.18Å 58.26Å 90.00° 100.72° 90.00°	Depositor
Resolution (Å)	30.00 – 1.80 30.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	91.8 (30.00-1.80) 91.9 (30.00-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 1.80Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.252 , 0.283 0.254 , 0.256	Depositor DCC
R_{free} test set	2978 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	21.3	Xtriage
Anisotropy	0.564	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 25.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.028 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5994	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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: OCS, PLP, EPE, TYO

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	0.79	5/2843 (0.2%)	1.17	21/3855 (0.5%)
2	B	0.76	4/3011 (0.1%)	1.14	21/4085 (0.5%)
All	All	0.78	9/5854 (0.2%)	1.16	42/7940 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	6
2	B	0	9
All	All	0	15

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	553	ILE	CA-CB	7.38	1.61	1.54
1	A	184	LEU	N-CA	6.97	1.55	1.46
2	B	724	GLN	N-CA	6.46	1.53	1.46
1	A	289	ARG	CA-C	6.25	1.60	1.52
1	A	160	ARG	N-CA	6.02	1.53	1.46
1	A	14	MET	C-N	-5.53	1.25	1.33
2	B	553	ILE	N-CA	5.28	1.53	1.46
1	A	291	ILE	CA-CB	5.24	1.60	1.54
2	B	602	ARG	N-CA	5.09	1.52	1.46

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	PRO	N-CA-C	12.64	132.06	111.34
1	A	184	LEU	N-CA-C	8.69	128.88	110.38
2	B	553	ILE	N-CA-CB	7.82	120.02	111.56
2	B	600	MET	CA-C-O	-7.80	112.47	120.66
1	A	53	ILE	N-CA-C	-6.74	98.03	107.80
2	B	635	ALA	N-CA-C	6.55	118.75	110.24
2	B	536	ASP	N-CA-C	6.51	119.37	111.82
1	A	248	THR	N-CA-C	-6.50	97.09	108.20
2	B	722	CYS	N-CA-C	6.46	119.92	110.17
2	B	748	THR	N-CA-C	-6.45	98.21	108.73
2	B	853	GLN	N-CA-C	6.25	118.09	111.28
1	A	340	GLU	N-CA-C	6.24	118.60	111.11
2	B	840	GLU	N-CA-C	6.21	118.60	111.02
2	B	725	VAL	N-CA-C	6.21	117.53	108.53
1	A	10	LEU	N-CA-C	6.18	119.31	110.10
1	A	182	VAL	N-CA-C	6.16	121.55	108.72
1	A	353	GLN	N-CA-C	6.08	117.98	111.36
1	A	321	HIS	N-CA-C	6.00	121.33	113.30
1	A	49	GLY	N-CA-C	-5.91	103.56	112.60
1	A	230	GLY	CA-C-N	5.88	125.50	119.56
1	A	230	GLY	C-N-CA	5.88	125.50	119.56
1	A	291	ILE	N-CA-CB	5.85	118.44	111.41
2	B	755	GLU	N-CA-C	5.80	117.88	109.07
2	B	681	PRO	N-CA-C	5.77	120.34	111.57
1	A	196	GLY	N-CA-C	-5.76	107.15	115.27
2	B	504	SER	N-CA-C	5.74	117.51	108.96
1	A	328	ARG	N-CA-C	5.68	117.23	110.19
2	B	809	PHE	N-CA-C	5.66	117.67	109.07
1	A	255	GLU	N-CA-C	5.64	117.65	109.07
2	B	606	ARG	N-CA-C	5.61	117.39	111.28
1	A	309	PHE	N-CA-C	5.46	117.36	109.07
1	A	107	LEU	CB-CA-C	-5.44	99.77	110.10
2	B	503	SER	N-CA-C	5.21	115.70	108.00
1	A	183[A]	SER	N-CA-C	5.18	121.83	110.80
1	A	183[B]	SER	N-CA-C	5.18	121.83	110.80
2	B	517	LYS	CA-C-N	5.18	125.16	120.03
2	B	517	LYS	C-N-CA	5.18	125.16	120.03
2	B	696	GLY	N-CA-C	-5.10	108.07	115.27
2	B	832	ILE	CA-C-N	5.05	126.15	119.84
2	B	832	ILE	C-N-CA	5.05	126.15	119.84
2	B	707	TYR	N-CA-C	5.04	119.42	113.12
1	A	287	VAL	N-CA-C	5.01	115.91	108.65

There are no chirality outliers.

All (15) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	158	PRO	Mainchain
1	A	182	VAL	Mainchain
1	A	289	ARG	Mainchain
2	B	551	PRO	Mainchain
2	B	600	MET	Mainchain
2	B	707	TYR	Sidechain
2	B	722	CYS	Mainchain
2	B	812	GLY	Mainchain

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	2799	0	2800	48	0
2	B	2944	0	2953	33	0
3	A	15	0	6	0	0
3	B	15	0	6	0	0
4	B	15	0	17	0	0
5	A	97	0	0	5	0
5	B	109	0	0	5	0
All	All	5994	0	5782	77	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 7.

All (77) 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:200:ASN:H	1:A:200:ASN:HD22	1.15	0.94
1:A:183[A]:SER:HB3	1:A:223:GLU:H	1.35	0.91
1:A:184:LEU:HD21	1:A:238:VAL:HG13	1.60	0.83
1:A:289:ARG:HG2	1:A:290[B]:THR:H	1.40	0.83
2:B:700:ASN:HD22	2:B:700:ASN:H	1.26	0.82
1:A:159[B]:ARG:O	1:A:160:ARG:HB2	1.78	0.82
1:A:290[A]:THR:O	5:A:402:HOH:O	1.98	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:683:SER:H	2:B:723[A]:GLU:HB2	1.55	0.70
1:A:73:GLN:HE22	2:B:571:SER:H	1.41	0.68
1:A:253:VAL:N	5:A:472:HOH:O	2.27	0.67
1:A:183[A]:SER:HB3	1:A:223:GLU:N	2.07	0.67
2:B:540:MET:HB3	2:B:552[B]:ARG:HB2	1.77	0.65
1:A:183[B]:SER:O	1:A:322:ARG:O	2.14	0.64
1:A:71:SER:H	2:B:573:GLN:HE22	1.46	0.63
1:A:200:ASN:H	1:A:200:ASN:ND2	1.93	0.62
1:A:200:ASN:HD22	1:A:200:ASN:N	1.95	0.62
2:B:813[B]:THR:O	2:B:814:ALA:HB2	1.99	0.62
2:B:723[A]:GLU:O	2:B:724:GLN:HB2	2.02	0.60
2:B:601[A]:LEU:HD13	2:B:611:SER:HB2	1.88	0.56
2:B:501:ALA:N	2:B:611:SER:HG	2.04	0.55
1:A:150:GLU:HG3	1:A:159[B]:ARG:HB2	1.89	0.54
2:B:834:THR:O	2:B:835:MET:HB2	2.07	0.54
2:B:601[A]:LEU:HA	2:B:604:ALA:HB3	1.89	0.53
1:A:5:PHE:HB2	1:A:51:PRO:HD3	1.90	0.53
1:A:43:TRP:HZ3	1:A:159[A]:ARG:HA	1.75	0.52
1:A:73:GLN:NE2	2:B:571:SER:H	2.06	0.52
2:B:607:LEU:HD11	2:B:703:LEU:HD22	1.92	0.52
1:A:290[A]:THR:HG22	5:A:447:HOH:O	2.10	0.51
2:B:715:GLN:HA	2:B:718:LEU:HD12	1.94	0.50
1:A:159[B]:ARG:O	1:A:160:ARG:CB	2.57	0.50
2:B:700:ASN:H	2:B:700:ASN:ND2	2.02	0.50
1:A:73:GLN:HE21	1:A:204:GLY:HA3	1.76	0.50
1:A:290[B]:THR:O	1:A:295:GLN:NE2	2.27	0.49
1:A:141:TYO:HE1	1:A:141:TYO:CZ	2.42	0.49
1:A:158:PRO:O	1:A:160:ARG:N	2.46	0.49
1:A:289:ARG:CG	1:A:290[B]:THR:H	2.10	0.49
1:A:328:ARG:HE	1:A:330:LEU:HD21	1.78	0.49
2:B:600:MET:O	2:B:602:ARG:N	2.46	0.49
2:B:813[B]:THR:HG23	5:B:69:HOH:O	2.12	0.49
2:B:573:GLN:HE21	2:B:704:GLY:HA3	1.78	0.48
2:B:597:MET:O	2:B:601[B]:LEU:HB2	2.14	0.48
2:B:813[B]:THR:O	2:B:814:ALA:CB	2.63	0.47
1:A:43:TRP:CZ3	1:A:159[A]:ARG:HA	2.50	0.46
1:A:71:SER:H	2:B:573:GLN:NE2	2.12	0.46
1:A:191:ILE:HD11	1:A:229:TYR:HD1	1.80	0.46
2:B:700:ASN:HD22	2:B:700:ASN:N	2.04	0.45
1:A:72:LEU:HD22	1:A:107:LEU:O	2.17	0.45
2:B:683:SER:OG	2:B:723[A]:GLU:HG2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:ILE:HD11	1:A:229:TYR:CD1	2.52	0.45
1:A:66:SER:HB2	1:A:72:LEU:HD12	1.99	0.45
2:B:552[A]:ARG:O	2:B:553:ILE:HB	2.17	0.45
2:B:706:ASN:ND2	5:B:112:HOH:O	2.51	0.44
1:A:292:THR:H	1:A:295:GLN:HE21	1.65	0.44
1:A:184:LEU:HD13	1:A:226:LEU:HB2	2.00	0.44
1:A:183[B]:SER:O	1:A:184:LEU:HB2	2.18	0.43
2:B:724:GLN:NE2	5:B:206:HOH:O	2.51	0.43
1:A:262:ASN:HD21	1:A:264:VAL:HG22	1.84	0.43
2:B:601[A]:LEU:O	2:B:605:MET:HB2	2.19	0.43
1:A:141:TYO:CZ	1:A:143:ARG:HH21	2.31	0.43
2:B:813[A]:THR:HG21	5:B:150:HOH:O	2.18	0.43
2:B:551:PRO:O	2:B:552[A]:ARG:HB3	2.18	0.43
1:A:290[B]:THR:O	5:A:464:HOH:O	2.21	0.42
1:A:46:LYS:H	1:A:46:LYS:HD2	1.84	0.42
2:B:517:LYS:HA	2:B:518:PRO:HD2	1.92	0.42
1:A:40:MET:HE2	1:A:54:GLN:HE22	1.84	0.42
1:A:258:THR:O	1:A:288:GLU:HA	2.19	0.42
1:A:183[B]:SER:HB2	1:A:322:ARG:N	2.34	0.41
1:A:140:LEU:HD21	1:A:167:LEU:HD13	2.02	0.41
1:A:322:ARG:HA	1:A:330:LEU:O	2.21	0.41
1:A:325:TYR:O	1:A:326:LYS:HG3	2.21	0.41
2:B:510:LEU:HD12	2:B:551:PRO:HB2	2.02	0.41
1:A:66:SER:HB3	5:A:450:HOH:O	2.20	0.41
1:A:318:CYS:HA	1:A:319:PRO:HD2	1.92	0.41
1:A:235:LEU:HD21	1:A:296:LEU:HD22	2.02	0.40
2:B:503:SER:HB3	2:B:504:SER:H	1.75	0.40
1:A:348:GLU:HG3	1:A:357:ARG:HH22	1.86	0.40
2:B:601[B]:LEU:HD22	5:B:17:HOH:O	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	345/365 (94%)	319 (92%)	16 (5%)	10 (3%)	3	0
2	B	366/365 (100%)	336 (92%)	18 (5%)	12 (3%)	3	0
All	All	711/730 (97%)	655 (92%)	34 (5%)	22 (3%)	5	0

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	45	ASP
1	A	159[A]	ARG
1	A	159[B]	ARG
1	A	160	ARG
1	A	172	ALA
1	A	183[A]	SER
1	A	183[B]	SER
1	A	290[A]	THR
1	A	290[B]	THR
2	B	601[A]	LEU
2	B	601[B]	LEU
2	B	602	ARG
2	B	723[A]	GLU
2	B	723[B]	GLU
2	B	813[A]	THR
2	B	813[B]	THR
2	B	636	ALA
2	B	814	ALA
2	B	835	MET
1	A	184	LEU
2	B	525	GLY
2	B	833	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	301/314 (96%)	280 (93%)	21 (7%)	14	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	B	318/315 (101%)	301 (95%)	17 (5%)	20 9
All	All	619/629 (98%)	581 (94%)	38 (6%)	18 6

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LEU
1	A	41	VAL
1	A	44	ASN
1	A	45	ASP
1	A	61	LEU
1	A	74	LEU
1	A	158	PRO
1	A	179	VAL
1	A	183[A]	SER
1	A	183[B]	SER
1	A	198	VAL
1	A	200	ASN
1	A	223	GLU
1	A	225	VAL
1	A	262	ASN
1	A	318	CYS
1	A	324	LEU
1	A	326	LYS
1	A	328	ARG
1	A	336	GLU
1	A	365	VAL
2	B	510	LEU
2	B	574	LEU
2	B	587	GLN
2	B	600	MET
2	B	601[A]	LEU
2	B	601[B]	LEU
2	B	607	LEU
2	B	619	GLU
2	B	698	VAL
2	B	700	ASN
2	B	796	LEU
2	B	813[A]	THR
2	B	813[B]	THR
2	B	818	CYS
2	B	835	MET

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Mol	Chain	Res	Type
2	B	846	GLN
2	B	865	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	44	ASN
1	A	54	GLN
1	A	57	GLN
1	A	73	GLN
1	A	157	GLN
1	A	200	ASN
1	A	206	ASN
1	A	249	HIS
1	A	262	ASN
1	A	295	GLN
2	B	511	GLN
2	B	554	GLN
2	B	573	GLN
2	B	586	GLN
2	B	587	GLN
2	B	657	GLN
2	B	700	ASN
2	B	706	ASN
2	B	724	GLN
2	B	734	GLN
2	B	772	GLN
2	B	821	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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
2	OCS	B	608	2	6,8,9	2.77	2 (33%)	7,11,13	1.79	2 (28%)
1	OCS	A	108	1	6,8,9	2.69	2 (33%)	7,11,13	2.28	3 (42%)
1	TYO	A	141	1	10,13,14	2.61	3 (30%)	3,14,16	2.00	1 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	OCS	B	608	2	-	1/4/7/9	-
1	OCS	A	108	1	-	1/4/7/9	-
1	TYO	A	141	1	-	5/11/14/16	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	108	OCS	CB-CA	-5.94	1.49	1.53
2	B	608	OCS	CB-CA	-5.82	1.49	1.53
1	A	141	TYO	CD1-CG	5.51	1.56	1.45
1	A	141	TYO	CE2-CD2	5.19	1.59	1.43
2	B	608	OCS	OD3-SG	-2.72	1.37	1.45
1	A	141	TYO	OE1-CE1	2.29	1.43	1.33
1	A	108	OCS	OD3-SG	-2.17	1.39	1.45

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	108	OCS	CA-CB-SG	4.10	119.70	113.61
1	A	141	TYO	CD1-CG-CD2	3.12	123.84	119.00
2	B	608	OCS	OD2-SG-OD3	2.88	118.59	111.40
1	A	108	OCS	OD2-SG-OD3	2.87	118.58	111.40
2	B	608	OCS	OD3-SG-CB	-2.82	102.54	106.76
1	A	108	OCS	OD1-SG-CB	-2.75	102.65	106.76

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	108	OCS	N-CA-CB-SG
1	A	141	TYO	CE1-CD1-CG-CB
1	A	141	TYO	CE2-CD2-CG-CB
1	A	141	TYO	CG-CD2-CE2-CZ
2	B	608	OCS	N-CA-CB-SG
1	A	141	TYO	CE2-CD2-CG-CD1
1	A	141	TYO	CE1-CD1-CG-CD2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	141	TYO	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	$\# Z > 2$
4	EPE	B	430	-	15,15,15	1.48	3 (20%)	19,20,20	0.90	2 (10%)
3	PLP	B	401	2	15,15,16	1.84	4 (26%)	21,22,23	1.85	7 (33%)
3	PLP	A	400	1	15,15,16	1.70	3 (20%)	21,22,23	1.72	6 (28%)

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
4	EPE	B	430	-	-	6/9/19/19	0/1/1/1
3	PLP	B	401	2	-	0/6/6/8	0/1/1/1
3	PLP	A	400	1	-	0/6/6/8	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	401	PLP	C2A-C2	4.18	1.57	1.50
3	A	400	PLP	C2A-C2	3.66	1.56	1.50
4	B	430	EPE	C10-S	3.00	1.81	1.77
3	A	400	PLP	C4A-C4	2.77	1.57	1.51
3	B	401	PLP	C4A-C4	2.72	1.57	1.51
4	B	430	EPE	C2-N1	2.61	1.53	1.46
4	B	430	EPE	C6-N1	2.57	1.53	1.46
3	A	400	PLP	C5-C4	2.17	1.43	1.40
3	B	401	PLP	C5-C4	2.12	1.42	1.40
3	B	401	PLP	O4P-C5A	-2.03	1.37	1.44

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	400	PLP	O4P-C5A-C5	3.68	116.25	109.36
3	B	401	PLP	O4P-C5A-C5	3.49	115.89	109.36
3	B	401	PLP	O2P-P-O4P	-3.19	98.36	106.67
3	A	400	PLP	O2P-P-O4P	-2.80	99.36	106.67
3	B	401	PLP	O4P-P-O1P	2.64	113.56	106.44
3	B	401	PLP	C5A-C5-C6	-2.61	115.10	119.36
3	B	401	PLP	O3P-P-O1P	2.58	120.90	110.83
3	B	401	PLP	C4A-C4-C5	2.46	123.48	120.94
3	A	400	PLP	C5A-C5-C6	-2.42	115.41	119.36
3	A	400	PLP	O3P-P-O1P	2.42	120.25	110.83
4	B	430	EPE	C9-N1-C6	-2.24	105.28	111.24
4	B	430	EPE	C7-N4-C3	2.21	117.13	111.24
3	A	400	PLP	C4A-C4-C5	2.18	123.18	120.94
3	A	400	PLP	O4P-P-O1P	2.11	112.14	106.44
3	B	401	PLP	C6-C5-C4	2.10	119.82	118.10

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	430	EPE	C9-C10-S-O1S

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Mol	Chain	Res	Type	Atoms
4	B	430	EPE	C9-C10-S-O2S
4	B	430	EPE	C9-C10-S-O3S
4	B	430	EPE	N4-C7-C8-O8
4	B	430	EPE	C10-C9-N1-C2
4	B	430	EPE	C10-C9-N1-C6

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	346/365 (94%)	1.12	60 (17%) 4 3	13, 22, 37, 47	3 (0%)
2	B	364/365 (99%)	1.16	61 (16%) 4 3	13, 22, 40, 48	4 (1%)
All	All	710/730 (97%)	1.14	121 (17%) 4 3	13, 22, 39, 48	7 (0%)

All (121) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	290[A]	THR	11.5
2	B	813[A]	THR	11.0
2	B	607	LEU	10.3
2	B	601[A]	LEU	8.9
2	B	723[A]	GLU	7.9
1	A	183[A]	SER	7.7
1	A	182	VAL	7.6
2	B	525	GLY	6.6
1	A	159[A]	ARG	6.5
2	B	551	PRO	6.4
2	B	511	GLN	6.3
1	A	175	PRO	6.1
2	B	526	GLU	5.8
2	B	552[A]	ARG	5.5
1	A	329	ASN	5.3
1	A	15	THR	5.2
2	B	815	CYS	5.2
1	A	184	LEU	5.1
1	A	328	ARG	5.1
2	B	677	GLY	5.1
1	A	315	CYS	4.9
2	B	812	GLY	4.8
2	B	501	ALA	4.7
2	B	718	LEU	4.6

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Mol	Chain	Res	Type	RSRZ
1	A	326	LYS	4.4
1	A	327	ASP	4.4
2	B	676	GLY	4.4
2	B	512	LEU	4.4
1	A	1	ALA	4.3
1	A	44	ASN	4.3
1	A	3	SER	4.2
1	A	338	GLY	4.2
2	B	517	LYS	4.1
2	B	502	SER	4.1
1	A	2	SER	4.0
2	B	722	CYS	3.9
2	B	527	PRO	3.9
1	A	150	GLU	3.8
1	A	107	LEU	3.8
1	A	10	LEU	3.8
2	B	600	MET	3.8
2	B	553	ILE	3.7
2	B	524	PRO	3.7
2	B	546	LYS	3.7
2	B	754	LEU	3.6
1	A	158	PRO	3.6
2	B	716	GLU	3.6
2	B	675	PRO	3.6
2	B	516	GLN	3.5
1	A	174	PHE	3.5
2	B	523	GLY	3.5
2	B	814	ALA	3.4
1	A	358	ALA	3.4
2	B	674	PHE	3.3
1	A	84	LYS	3.2
1	A	87	GLN	3.2
1	A	194	TRP	3.1
2	B	678	SER	3.1
1	A	83	GLY	3.1
2	B	694	TRP	3.1
1	A	171	GLY	3.1
2	B	827	ASP	3.0
1	A	176	GLY	3.0
2	B	602	ARG	3.0
2	B	528	LEU	2.9
1	A	331	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	356	ILE	2.9
1	A	223	GLU	2.9
1	A	134	ASP	2.9
2	B	510	LEU	2.9
1	A	33	THR	2.8
2	B	636	ALA	2.8
1	A	323	ILE	2.8
1	A	330	LEU	2.8
1	A	160	ARG	2.8
2	B	779	GLN	2.8
1	A	198	VAL	2.7
2	B	506	LYS	2.7
1	A	172	ALA	2.6
1	A	173	TYR	2.6
2	B	630	ASP	2.6
1	A	12	LEU	2.6
2	B	545	ASP	2.6
2	B	834	THR	2.6
2	B	530	PHE	2.6
1	A	178	SER	2.6
2	B	503	SER	2.6
1	A	359	HIS	2.5
1	A	133	PRO	2.5
1	A	43	TRP	2.5
1	A	318	CYS	2.5
1	A	325	TYR	2.5
2	B	865	VAL	2.5
1	A	46	LYS	2.5
1	A	85	ASP	2.5
2	B	529	VAL	2.4
2	B	679	VAL	2.4
1	A	177	GLY	2.4
2	B	611	SER	2.4
2	B	672	ALA	2.4
1	A	52	ARG	2.4
2	B	533	THR	2.3
1	A	289	ARG	2.3
2	B	691	ILE	2.3
2	B	857	ARG	2.3
1	A	179	VAL	2.3
1	A	350	LYS	2.3
2	B	604	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	520	LYS	2.2
1	A	61	LEU	2.2
1	A	218	LEU	2.2
1	A	80	ALA	2.2
2	B	515	THR	2.1
2	B	514	MET	2.1
1	A	365	VAL	2.1
1	A	191	ILE	2.1
2	B	790	THR	2.1
2	B	801	GLU	2.1
1	A	225	VAL	2.1
2	B	585	ASP	2.1
2	B	605	MET	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	OCS	A	108	9/10	0.66	0.27	29,33,39,39	0
2	OCS	B	608	9/10	0.69	0.21	28,32,38,38	0
1	TYO	A	141	14/15	0.77	0.17	19,25,34,37	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(Å ²)	Q<0.9
4	EPE	B	430	15/15	0.74	0.18	37,40,46,46	0
3	PLP	A	400	15/16	0.95	0.09	17,19,20,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PLP	B	401	15/16	0.96	0.07	15,19,22,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.