



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:42 PM GMT

PDB ID : 1Q8A
Title : Cobalamin-dependent methionine synthase (1-566) from *Thermotoga maritima* (Cd2+:L-Hcy complex, Se-Met)
Authors : Evans, J.C.; Huddler, D.P.; Hilgers, M.T.; Romanchuk, G.; Matthews, R.G.; Ludwig, M.L.
Deposited on : 2003-08-20
Resolution : 1.70 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

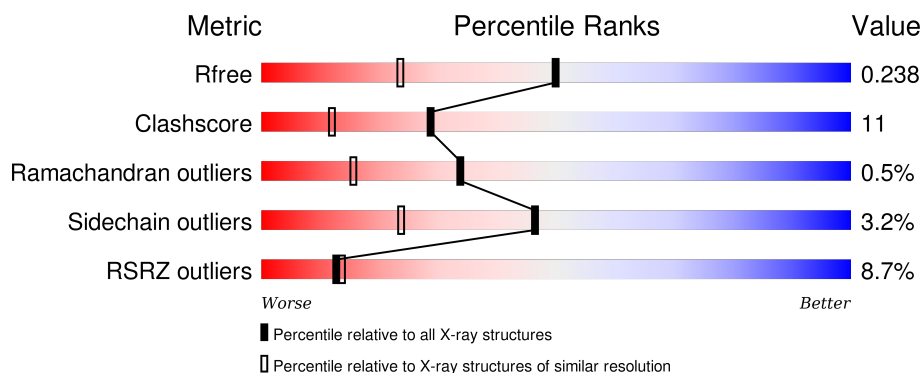
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	91344	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	566	<div> <div>8%</div> <div>76%</div> <div>21%</div> <div>..</div> </div>
1	B	566	<div> <div>9%</div> <div>77%</div> <div>17%</div> <div>..</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9007 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 5-methyltetrahydrofolate S-homocysteine methyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	559	Total	C	N	O	S	Se	0	0	0
			4423	2839	738	833	3	10			
1	B	547	Total	C	N	O	S	Se	0	0	0
			4333	2784	722	814	3	10			

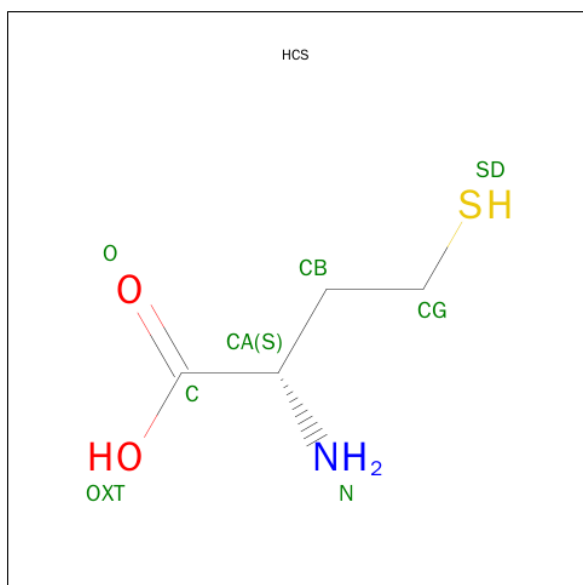
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
A	27	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
A	71	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
A	135	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
A	174	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
A	334	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
A	422	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
A	439	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
A	537	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
A	545	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	27	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	71	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	135	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	174	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	334	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	422	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	439	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	537	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5
B	545	MSE	MET	MODIFIED RESIDUE	UNP Q9WYA5

- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Cd 1 1	0	0
2	A	1	Total Cd 1 1	0	0

- Molecule 3 is 2-AMINO-4-MERCAPTO-BUTYRIC ACID (three-letter code: HCS) (formula: C₄H₉NO₂S).



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

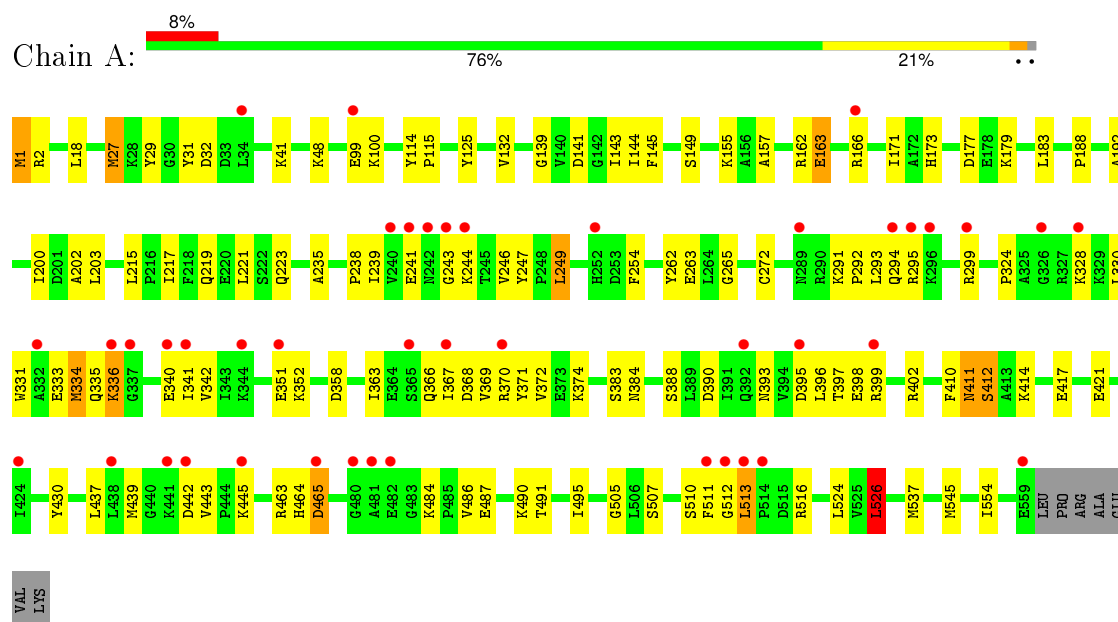
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	135	Total O 135 135	0	0
4	B	98	Total O 98 98	0	0

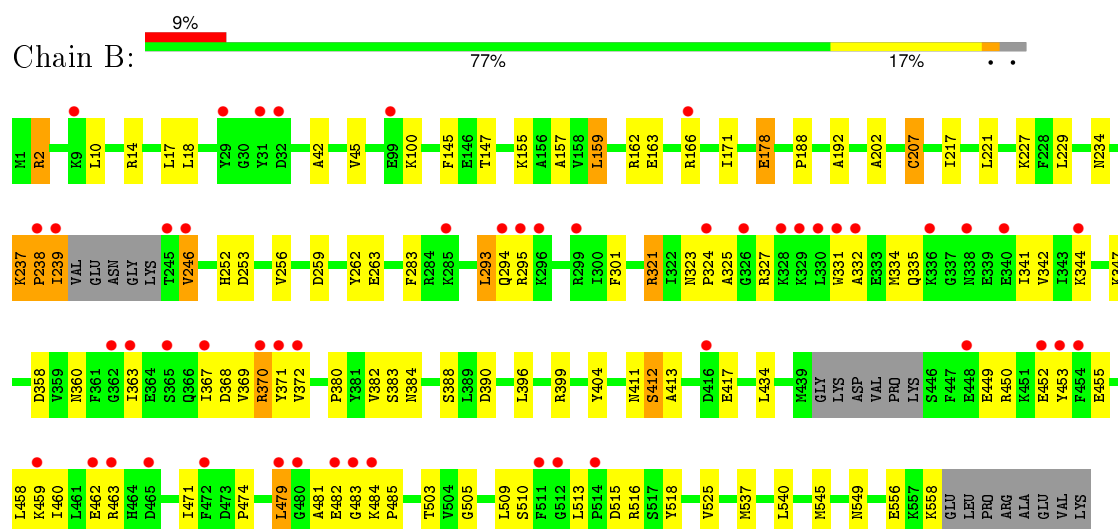
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: 5-methyltetrahydrofolate S-homocysteine methyltransferase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.12Å 86.19Å 126.08Å 90.00° 99.99° 90.00°	Depositor
Resolution (Å)	20.00 – 1.70 17.92 – 1.65	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-1.70) 90.4 (17.92-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 1.65Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.214 , 0.242 0.210 , 0.238	Depositor DCC
R_{free} test set	12710 reflections (10.10%)	DCC
Wilson B-factor (Å ²)	21.5	Xtriage
Anisotropy	0.399	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 62.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 134815 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9007	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: HCS, CD

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	0.66	2/4496 (0.0%)	0.83	5/6056 (0.1%)
1	B	0.55	0/4403	0.76	3/5929 (0.1%)
All	All	0.61	2/8899 (0.0%)	0.80	8/11985 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	334	MSE	SE-CE	-5.39	1.63	1.95
1	A	272	CYS	CB-SG	5.08	1.90	1.82

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	412	SER	N-CA-C	9.04	135.42	111.00
1	B	412	SER	N-CA-C	6.68	129.03	111.00
1	A	411	ASN	C-N-CA	6.15	137.07	121.70
1	B	505	GLY	N-CA-C	-5.71	98.83	113.10
1	A	411	ASN	CA-C-N	-5.59	104.91	117.20
1	A	505	GLY	N-CA-C	-5.52	99.30	113.10
1	B	321	ARG	N-CA-C	5.28	125.24	111.00
1	A	526	LEU	CA-CB-CG	5.08	126.98	115.30

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	4423	0	4510	121	0
1	B	4333	0	4416	85	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	8	0	7	0	0
3	B	8	0	7	0	0
4	A	135	0	0	5	0
4	B	98	0	0	0	0
All	All	9007	0	8940	203	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 11.

All (203) 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:372:VAL:HG21	1:B:396:LEU:HD11	1.20	1.15
1:A:241:GLU:HG2	1:A:246:VAL:HG11	1.40	1.02
1:A:331:TRP:CH2	1:A:367:ILE:HD11	2.11	0.86
1:A:27:MSE:HE3	1:A:32:ASP:N	1.94	0.81
1:A:241:GLU:HG2	1:A:246:VAL:CG1	2.10	0.79
1:A:330:LEU:O	1:A:334:MSE:HG3	1.83	0.79
1:B:321:ARG:NE	1:B:540:LEU:HD22	1.97	0.78
1:B:449:GLU:O	1:B:452:GLU:HG2	1.83	0.77
1:A:328:LYS:N	1:A:328:LYS:HE2	1.99	0.77
1:A:369:VAL:CG2	1:A:399:ARG:HD3	2.14	0.77
1:B:458:LEU:O	1:B:462:GLU:HG2	1.84	0.76
1:A:241:GLU:CG	1:A:246:VAL:HG11	2.16	0.75
1:A:331:TRP:HH2	1:A:367:ILE:HD11	1.50	0.75
1:A:484:LYS:HG2	1:A:487:GLU:HG2	1.69	0.74
1:A:369:VAL:HG21	1:A:399:ARG:HD3	1.67	0.74
1:A:486:VAL:HG12	1:A:490:LYS:HE3	1.71	0.71
1:A:484:LYS:HD2	1:B:556:GLU:OE2	1.90	0.71
1:A:336:LYS:NZ	1:A:336:LYS:HB3	2.06	0.71
1:A:363:ILE:H	1:A:363:ILE:HD12	1.57	0.70
1:A:367:ILE:N	1:A:367:ILE:HD12	2.06	0.70
1:A:333:GLU:OE2	1:A:336:LYS:HE2	1.92	0.70
1:A:363:ILE:N	1:A:363:ILE:HD12	2.08	0.69
1:A:340:GLU:HB2	4:A:828:HOH:O	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:372:VAL:HG21	1:B:396:LEU:CD1	2.13	0.68
1:B:390:ASP:HA	1:B:411:ASN:HB3	1.73	0.68
1:A:100:LYS:HE3	4:A:836:HOH:O	1.93	0.67
1:B:239:ILE:HD13	1:B:239:ILE:C	2.15	0.66
1:B:331:TRP:O	1:B:335:GLN:HG2	1.96	0.65
1:B:324:PRO:HG3	1:B:334:MSE:SE	2.46	0.65
1:A:484:LYS:CG	1:A:487:GLU:HG2	2.27	0.65
1:A:369:VAL:HG21	1:A:399:ARG:CD	2.28	0.64
1:B:227:LYS:O	1:B:295:ARG:NH2	2.29	0.64
1:B:369:VAL:O	1:B:372:VAL:HG22	1.98	0.64
1:B:159:LEU:O	1:B:163:GLU:HG3	1.97	0.64
1:A:510:SER:HA	1:A:513:LEU:HD21	1.78	0.64
1:A:192:ALA:HB2	1:A:221:LEU:HD12	1.80	0.64
1:B:341:ILE:O	1:B:344:LYS:HB3	1.98	0.64
1:A:1:MSE:SE	1:A:139:GLY:O	2.66	0.64
1:B:363:ILE:H	1:B:363:ILE:HD12	1.63	0.63
1:B:259:ASP:O	1:B:263:GLU:HG2	1.98	0.63
1:A:299:ARG:HH21	1:A:299:ARG:HG2	1.63	0.62
1:A:238:PRO:HG3	1:A:247:TYR:HE1	1.65	0.62
1:B:347:LYS:HE3	1:B:382:VAL:HG22	1.82	0.62
1:B:331:TRP:CH2	1:B:367:ILE:HD11	2.34	0.61
1:A:336:LYS:HZ3	1:A:336:LYS:HB3	1.65	0.61
1:A:166:ARG:HD2	1:A:166:ARG:N	2.15	0.61
1:A:114:TYR:CZ	1:A:374:LYS:HE2	2.36	0.60
1:A:1:MSE:SE	1:A:1:MSE:N	2.85	0.60
1:A:397:THR:HG23	1:A:410:PHE:CE1	2.37	0.60
1:B:331:TRP:CZ3	1:B:367:ILE:HD11	2.38	0.59
1:B:481:ALA:O	1:B:482:GLU:HB2	2.01	0.59
1:A:513:LEU:N	1:A:513:LEU:HD23	2.18	0.59
1:B:332:ALA:O	1:B:335:GLN:HB2	2.03	0.58
1:A:238:PRO:HG3	1:A:247:TYR:CE1	2.38	0.58
1:A:200:ILE:HD11	1:A:203:LEU:HD21	1.84	0.58
1:A:414:LYS:NZ	1:A:439:MSE:SE	2.86	0.58
1:B:325:ALA:O	1:B:327:ARG:HG3	2.04	0.57
1:A:328:LYS:CA	1:A:328:LYS:HE2	2.34	0.57
1:A:486:VAL:O	1:A:490:LYS:HG3	2.04	0.57
1:A:366:GLN:C	1:A:367:ILE:HD12	2.25	0.57
1:A:1:MSE:SE	1:A:1:MSE:H1	2.38	0.57
1:A:239:ILE:HG13	1:A:246:VAL:HG13	1.85	0.56
1:A:484:LYS:HG2	1:A:487:GLU:CG	2.35	0.56
1:A:200:ILE:CD1	1:A:203:LEU:HD21	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:479:LEU:C	1:B:481:ALA:H	2.08	0.56
1:B:166:ARG:HG3	1:B:166:ARG:HH11	1.71	0.56
1:A:294:GLN:HA	1:A:294:GLN:OE1	2.06	0.56
1:B:10:LEU:HD11	1:B:14:ARG:NH1	2.21	0.56
1:A:390:ASP:HA	1:A:411:ASN:HB3	1.89	0.55
1:A:166:ARG:HD3	4:A:767:HOH:O	2.07	0.55
1:B:549:ASN:OD1	1:B:558:LYS:NZ	2.36	0.55
1:A:188:PRO:HG3	1:A:217:ILE:HG23	1.89	0.54
1:B:42:ALA:O	1:B:45:VAL:HG12	2.07	0.54
1:B:178:GLU:H	1:B:178:GLU:CD	2.11	0.54
1:A:331:TRP:CZ3	1:A:367:ILE:HD11	2.41	0.54
1:B:229:LEU:HG	1:B:295:ARG:NH2	2.22	0.54
1:A:443:VAL:CG2	1:A:445:LYS:HE2	2.38	0.54
1:B:367:ILE:N	1:B:367:ILE:HD12	2.22	0.54
1:A:27:MSE:HE3	1:A:32:ASP:HB3	1.90	0.54
1:B:207:CYS:HA	1:B:234:ASN:OD1	2.08	0.54
1:B:460:ILE:HA	1:B:463:ARG:HH11	1.73	0.53
1:B:510:SER:OG	1:B:516:ARG:HB2	2.09	0.53
1:A:99:GLU:OE1	1:A:99:GLU:N	2.32	0.53
1:B:321:ARG:CZ	1:B:540:LEU:HD22	2.39	0.53
1:A:239:ILE:O	1:A:246:VAL:HG12	2.09	0.52
1:A:331:TRP:CZ2	1:A:363:ILE:HD13	2.44	0.52
1:A:336:LYS:NZ	1:A:336:LYS:CB	2.68	0.52
1:A:27:MSE:HE3	1:A:32:ASP:CB	2.40	0.52
1:B:358:ASP:HA	1:B:388:SER:HB3	1.92	0.52
1:A:200:ILE:CD1	1:A:203:LEU:CD2	2.89	0.51
1:A:398:GLU:O	1:A:402:ARG:HG2	2.11	0.51
1:B:341:ILE:HG23	1:B:342:VAL:N	2.24	0.51
1:A:177:ASP:OD1	1:A:179:LYS:HG3	2.10	0.51
1:B:515:ASP:HB3	1:B:518:TYR:HD1	1.76	0.51
1:B:479:LEU:HD21	1:B:513:LEU:CD1	2.40	0.51
1:B:238:PRO:O	1:B:246:VAL:O	2.29	0.50
1:A:155:LYS:HE2	1:A:402:ARG:O	2.11	0.50
1:B:479:LEU:N	1:B:479:LEU:HD23	2.27	0.50
1:B:331:TRP:CZ2	1:B:363:ILE:HD13	2.47	0.49
1:A:363:ILE:CD1	1:A:363:ILE:H	2.24	0.49
1:A:27:MSE:SE	1:A:32:ASP:HB3	2.61	0.49
1:A:486:VAL:CG1	1:A:490:LYS:HE3	2.39	0.49
1:B:166:ARG:NH1	1:B:166:ARG:HG3	2.27	0.49
1:B:537:MSE:HE2	1:B:545:MSE:SE	2.62	0.49
1:A:299:ARG:HG2	1:A:299:ARG:NH2	2.26	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:GLU:OE2	1:A:352:LYS:HD2	2.12	0.49
1:A:27:MSE:HE3	1:A:31:TYR:C	2.32	0.49
1:A:324:PRO:HA	1:A:330:LEU:HD23	1.94	0.49
1:A:1:MSE:HG2	1:A:141:ASP:OD2	2.12	0.49
1:B:479:LEU:HD22	1:B:509:LEU:O	2.12	0.49
1:B:145:PHE:CE2	1:B:157:ALA:HB1	2.48	0.49
1:B:479:LEU:CD2	1:B:479:LEU:H	2.25	0.49
1:A:328:LYS:HA	1:A:328:LYS:HE2	1.95	0.48
1:B:363:ILE:N	1:B:363:ILE:HD12	2.29	0.48
1:A:188:PRO:CG	1:A:217:ILE:HG23	2.43	0.48
1:B:450:ARG:NH2	1:B:474:PRO:O	2.47	0.48
1:B:479:LEU:HD21	1:B:513:LEU:HD11	1.96	0.48
1:B:207:CYS:SG	1:B:234:ASN:OD1	2.72	0.48
1:A:383:SER:O	1:A:384:ASN:HB2	2.14	0.48
1:B:332:ALA:HA	1:B:335:GLN:HG3	1.96	0.47
1:A:437:LEU:HG	1:A:439:MSE:HG2	1.96	0.47
1:B:253:ASP:O	1:B:256:VAL:HG22	2.14	0.47
1:B:474:PRO:HD2	1:B:503:THR:O	2.14	0.47
1:B:301:PHE:CE1	1:B:471:ILE:HD11	2.49	0.47
1:A:219:GLN:O	1:A:223:GLN:HG3	2.14	0.47
1:A:507:SER:HB3	1:A:511:PHE:CE1	2.50	0.47
1:B:479:LEU:CD2	1:B:509:LEU:O	2.63	0.47
1:A:443:VAL:HG21	1:A:445:LYS:HE2	1.96	0.47
1:B:301:PHE:CZ	1:B:471:ILE:HG13	2.50	0.47
1:B:155:LYS:HE3	1:B:404:TYR:O	2.14	0.47
1:A:29:TYR:HE2	1:A:48:LYS:HE3	1.79	0.47
1:B:2:ARG:HD3	1:B:2:ARG:N	2.30	0.47
1:A:333:GLU:HA	1:A:336:LYS:HG2	1.96	0.47
1:A:27:MSE:HA	1:A:27:MSE:HE2	1.97	0.47
1:A:265:GLY:HA3	1:A:295:ARG:HD3	1.96	0.47
1:B:188:PRO:HG2	1:B:217:ILE:HG23	1.97	0.47
1:A:239:ILE:CG1	1:A:246:VAL:HG13	2.45	0.47
1:A:491:THR:O	1:A:495:ILE:HG12	2.15	0.46
1:A:372:VAL:HG21	1:A:396:LEU:HD21	1.97	0.46
1:A:125:TYR:OH	1:A:163:GLU:HG3	2.15	0.46
1:B:188:PRO:CG	1:B:217:ILE:HG23	2.45	0.46
1:A:341:ILE:HG23	1:A:342:VAL:N	2.30	0.46
1:A:145:PHE:CE2	1:A:157:ALA:HB1	2.51	0.46
1:B:368:ASP:OD1	1:B:370:ARG:HG2	2.16	0.46
1:B:239:ILE:CD1	1:B:239:ILE:C	2.83	0.45
1:A:554:ILE:HG23	1:B:485:PRO:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:LYS:HD3	1:A:244:LYS:HA	1.79	0.45
1:A:27:MSE:HE3	1:A:32:ASP:CA	2.46	0.45
1:A:291:LYS:HG3	1:A:292:PRO:HD2	1.97	0.45
1:A:367:ILE:N	1:A:367:ILE:CD1	2.77	0.45
1:A:166:ARG:CD	4:A:767:HOH:O	2.65	0.45
1:B:171:ILE:HG12	1:B:202:ALA:HB3	1.98	0.45
1:B:417:GLU:HG2	1:B:460:ILE:CD1	2.47	0.44
1:B:262:TYR:CZ	1:B:293:LEU:HD13	2.52	0.44
1:A:41:LYS:HB2	1:A:41:LYS:HE3	1.84	0.44
1:A:537:MSE:HE2	1:A:545:MSE:SE	2.67	0.44
1:A:2:ARG:NH2	1:A:99:GLU:O	2.50	0.44
1:A:369:VAL:HG13	1:A:370:ARG:N	2.33	0.44
1:A:398:GLU:HG3	1:A:430:TYR:CE2	2.53	0.44
1:B:239:ILE:O	1:B:239:ILE:HG23	2.18	0.44
1:A:393:ASN:OD1	1:A:395:ASP:HB3	2.18	0.44
1:A:235:ALA:C	1:A:249:LEU:HB2	2.39	0.44
1:B:479:LEU:CD2	1:B:479:LEU:N	2.80	0.43
1:A:331:TRP:O	1:A:335:GLN:HG3	2.18	0.43
1:B:192:ALA:HB2	1:B:221:LEU:HD12	1.99	0.43
1:A:336:LYS:HZ2	1:A:336:LYS:HB3	1.83	0.43
1:A:239:ILE:HG13	1:A:246:VAL:CG1	2.48	0.43
1:A:27:MSE:HE2	1:A:31:TYR:N	2.34	0.43
1:A:368:ASP:O	1:A:371:TYR:N	2.50	0.43
1:B:479:LEU:HA	1:B:483:GLY:HA3	2.00	0.43
1:B:484:LYS:N	1:B:485:PRO:HD3	2.34	0.42
1:B:162:ARG:HA	1:B:162:ARG:HD3	1.69	0.42
1:A:144:ILE:HD11	1:A:173:HIS:CE1	2.55	0.42
1:B:178:GLU:N	1:B:178:GLU:CD	2.72	0.42
1:A:417:GLU:OE1	1:A:463:ARG:NH1	2.48	0.42
1:B:323:ASN:HA	1:B:324:PRO:HD3	1.85	0.42
1:A:510:SER:OG	1:A:516:ARG:HB2	2.20	0.42
1:A:262:TYR:CZ	1:A:293:LEU:HD13	2.55	0.42
1:B:360:ASN:HB2	1:B:390:ASP:HB3	2.02	0.42
1:A:507:SER:HB3	1:A:511:PHE:CZ	2.55	0.42
4:A:826:HOH:O	1:B:485:PRO:HD2	2.18	0.42
1:A:263:GLU:HA	1:A:293:LEU:CD2	2.50	0.42
1:A:526:LEU:HD13	1:B:525:VAL:HG12	2.02	0.42
1:A:243:GLY:O	1:A:244:LYS:HG2	2.20	0.41
1:A:368:ASP:O	1:A:371:TYR:HB3	2.21	0.41
1:A:149:SER:HB3	1:A:183:LEU:HB3	2.03	0.41
1:A:358:ASP:HA	1:A:388:SER:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:516:ARG:HG2	1:A:516:ARG:HH11	1.86	0.41
1:B:237:LYS:HG3	1:B:238:PRO:HD2	2.02	0.41
1:B:370:ARG:H	1:B:370:ARG:HG2	1.59	0.41
1:B:413:ALA:HB2	1:B:434:LEU:HD11	2.02	0.41
1:A:464:HIS:O	1:A:465:ASP:C	2.59	0.41
1:A:132:VAL:HG13	1:A:143:ILE:CD1	2.51	0.41
1:B:383:SER:O	1:B:384:ASN:HB2	2.21	0.41
1:A:162:ARG:HA	1:A:162:ARG:HD3	1.79	0.41
1:B:459:LYS:O	1:B:463:ARG:HG3	2.21	0.40
1:B:368:ASP:O	1:B:371:TYR:N	2.46	0.40
1:A:417:GLU:O	1:A:421:GLU:HG3	2.21	0.40
1:B:17:LEU:HD11	1:B:283:PHE:HB3	2.04	0.40
1:A:171:ILE:HG12	1:A:202:ALA:HB3	2.03	0.40
1:A:166:ARG:HD2	1:A:166:ARG:H	1.83	0.40
1:A:114:TYR:CE2	1:A:374:LYS:HE2	2.56	0.40
1:B:380:PRO:O	1:B:384:ASN:HA	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	557/566 (98%)	531 (95%)	23 (4%)	3 (0%)	34	15
1	B	541/566 (96%)	515 (95%)	24 (4%)	2 (0%)	39	20
All	All	1098/1132 (97%)	1046 (95%)	47 (4%)	5 (0%)	34	15

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	412	SER
1	A	512	GLY

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Mol	Chain	Res	Type
1	B	238	PRO
1	B	412	SER
1	A	465	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	488/484 (101%)	475 (97%)	13 (3%)	52	31
1	B	478/484 (99%)	460 (96%)	18 (4%)	40	17
All	All	966/968 (100%)	935 (97%)	31 (3%)	46	24

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	18	LEU
1	A	27	MSE
1	A	115	PRO
1	A	163	GLU
1	A	215	LEU
1	A	249	LEU
1	A	254	PHE
1	A	336	LYS
1	A	442	ASP
1	A	513	LEU
1	A	524	LEU
1	A	526	LEU
1	B	2	ARG
1	B	18	LEU
1	B	100	LYS
1	B	147	THR
1	B	159	LEU
1	B	178	GLU
1	B	207	CYS

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Mol	Chain	Res	Type
1	B	237	LYS
1	B	239	ILE
1	B	246	VAL
1	B	252	HIS
1	B	293	LEU
1	B	294	GLN
1	B	370	ARG
1	B	399	ARG
1	B	453	TYR
1	B	455	GLU
1	B	479	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	HCS	A	701	-	3,7,7	0.29	0	2,8,8	0.63	0
3	HCS	B	702	-	3,7,7	0.69	0	2,8,8	0.56	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	HCS	A	701	-	-	0/3/7/7	0/0/0/0
3	HCS	B	702	-	-	0/3/7/7	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	549/566 (96%)	0.37	43 (7%) 16 18	14, 25, 46, 59	0
1	B	537/566 (94%)	0.61	51 (9%) 10 11	17, 31, 52, 58	0
All	All	1086/1132 (95%)	0.49	94 (8%) 13 14	14, 28, 50, 59	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	239	ILE	11.5
1	A	511	PHE	11.3
1	B	245	THR	8.2
1	A	242	ASN	7.3
1	A	243	GLY	7.2
1	B	363	ILE	6.1
1	B	511	PHE	5.8
1	B	328	LYS	5.7
1	B	296	LYS	5.1
1	A	328	LYS	5.0
1	B	238	PRO	4.9
1	A	296	LYS	4.8
1	B	512	GLY	4.7
1	A	244	LYS	4.6
1	B	482	GLU	4.6
1	B	32	ASP	4.5
1	B	246	VAL	4.5
1	A	480	GLY	4.4
1	A	512	GLY	4.4
1	B	463	ARG	4.1
1	A	514	PRO	4.0
1	A	241	GLU	4.0
1	A	441	LYS	4.0
1	B	340	GLU	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	240	VAL	3.9
1	B	465	ASP	3.8
1	A	336	LYS	3.8
1	A	442	ASP	3.7
1	B	29	TYR	3.6
1	B	367	ILE	3.6
1	B	370	ARG	3.5
1	B	294	GLN	3.5
1	B	326	GLY	3.5
1	A	326	GLY	3.4
1	A	365	SER	3.4
1	B	514	PRO	3.4
1	B	336	LYS	3.4
1	B	331	TRP	3.4
1	A	513	LEU	3.3
1	A	294	GLN	3.3
1	A	465	ASP	3.2
1	B	483	GLY	3.1
1	B	479	LEU	3.0
1	B	365	SER	3.0
1	B	484	LYS	3.0
1	A	481	ALA	2.9
1	B	453	TYR	2.8
1	A	99	GLU	2.8
1	A	482	GLU	2.8
1	B	448	GLU	2.7
1	A	445	LYS	2.7
1	A	424	ILE	2.7
1	A	252	HIS	2.7
1	A	399	ARG	2.7
1	B	299	ARG	2.7
1	A	299	ARG	2.7
1	A	367	ILE	2.6
1	B	295	ARG	2.6
1	A	392	GLN	2.6
1	A	351	GLU	2.6
1	B	462	GLU	2.5
1	B	480	GLY	2.5
1	B	166	ARG	2.5
1	A	395	ASP	2.5
1	B	452	GLU	2.5
1	B	329	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	295	ARG	2.4
1	B	332	ALA	2.4
1	B	285	LYS	2.3
1	A	337	GLY	2.3
1	A	332	ALA	2.3
1	B	324	PRO	2.3
1	B	416	ASP	2.3
1	A	438	LEU	2.2
1	B	330	LEU	2.2
1	B	344	LYS	2.2
1	B	31	TYR	2.2
1	B	371	TYR	2.2
1	B	362	GLY	2.2
1	B	99	GLU	2.2
1	A	344	LYS	2.1
1	A	289	ASN	2.1
1	A	340	GLU	2.1
1	A	34	LEU	2.1
1	A	370	ARG	2.1
1	A	559	GLU	2.1
1	A	341	ILE	2.1
1	B	372	VAL	2.0
1	A	166	ARG	2.0
1	B	338	ASN	2.0
1	B	454	PHE	2.0
1	B	472	PHE	2.0
1	B	9	LYS	2.0
1	B	459	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	HCS	A	701	8/8	0.98	0.07	-0.36	14,19,23,23	0
3	HCS	B	702	8/8	0.99	0.05	-0.96	15,21,22,23	0
2	CD	B	602	1/1	1.00	0.04	-1.01	23,23,23,23	0
2	CD	A	601	1/1	0.99	0.03	-4.06	21,21,21,21	0

6.5 Other polymers

There are no such residues in this entry.