



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:47 PM UTC

PDB ID : 3KHK / pdb_00003khh
Title : Crystal structure of type-I restriction-modification system methylation subunit (MM_0429) from Methanosarchina mazei.
Authors : Ramagopal, U.A.; Toro, R.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-10-30
Resolution : 2.55 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

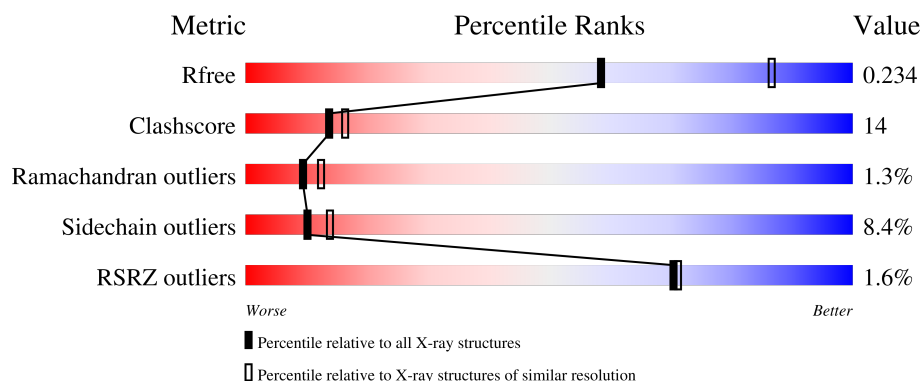
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

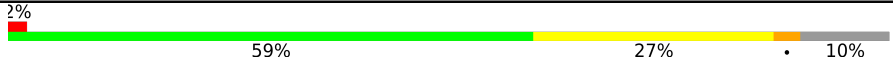
The reported resolution of this entry is 2.55 Å.

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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	544	
1	B	544	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7891 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 Type I restriction-modification system methylation subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	0	0
			3934	2514	669	736	15			
1	B	487	Total	C	N	O	S	0	0	0
			3929	2509	666	739	15			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP Q8PZR3
A	1	SER	-	expression tag	UNP Q8PZR3
A	2	LEU	-	expression tag	UNP Q8PZR3
A	536	GLU	-	expression tag	UNP Q8PZR3
A	537	GLY	-	expression tag	UNP Q8PZR3
A	538	HIS	-	expression tag	UNP Q8PZR3
A	539	HIS	-	expression tag	UNP Q8PZR3
A	540	HIS	-	expression tag	UNP Q8PZR3
A	541	HIS	-	expression tag	UNP Q8PZR3
A	542	HIS	-	expression tag	UNP Q8PZR3
A	543	HIS	-	expression tag	UNP Q8PZR3
B	0	MET	-	expression tag	UNP Q8PZR3
B	1	SER	-	expression tag	UNP Q8PZR3
B	2	LEU	-	expression tag	UNP Q8PZR3
B	536	GLU	-	expression tag	UNP Q8PZR3
B	537	GLY	-	expression tag	UNP Q8PZR3
B	538	HIS	-	expression tag	UNP Q8PZR3
B	539	HIS	-	expression tag	UNP Q8PZR3
B	540	HIS	-	expression tag	UNP Q8PZR3
B	541	HIS	-	expression tag	UNP Q8PZR3
B	542	HIS	-	expression tag	UNP Q8PZR3
B	543	HIS	-	expression tag	UNP Q8PZR3

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

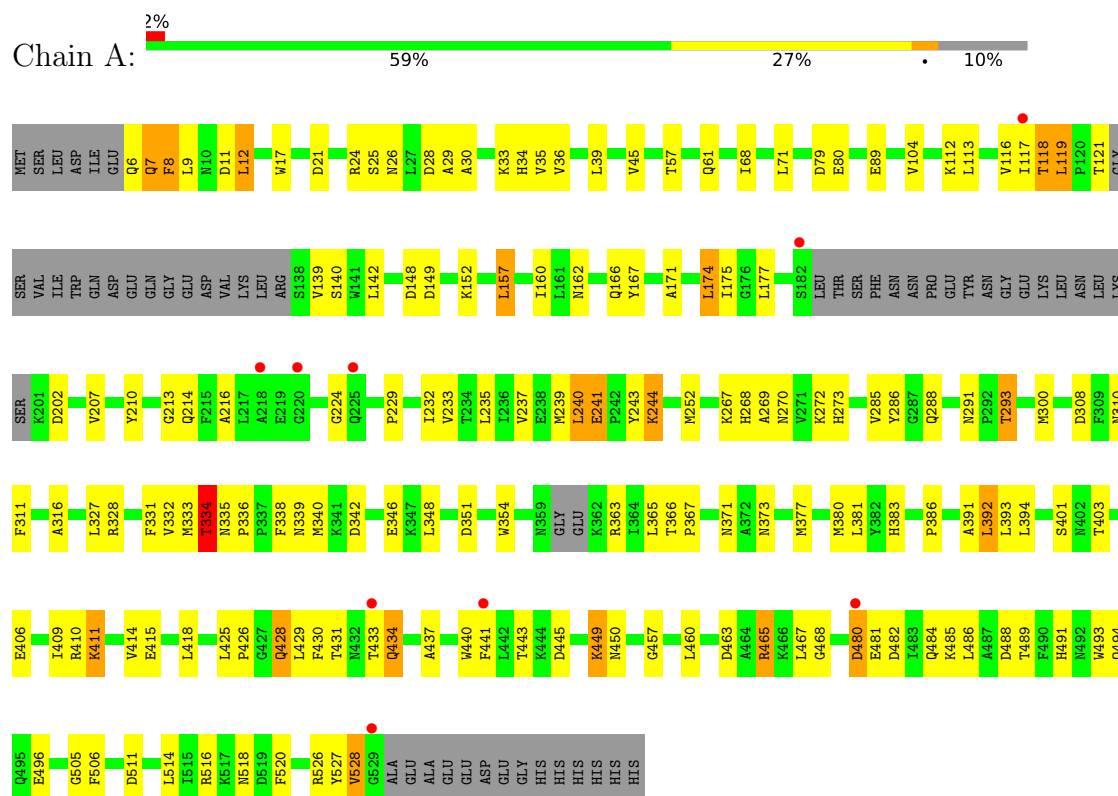
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	7	Total	O	0	0
			7	7		

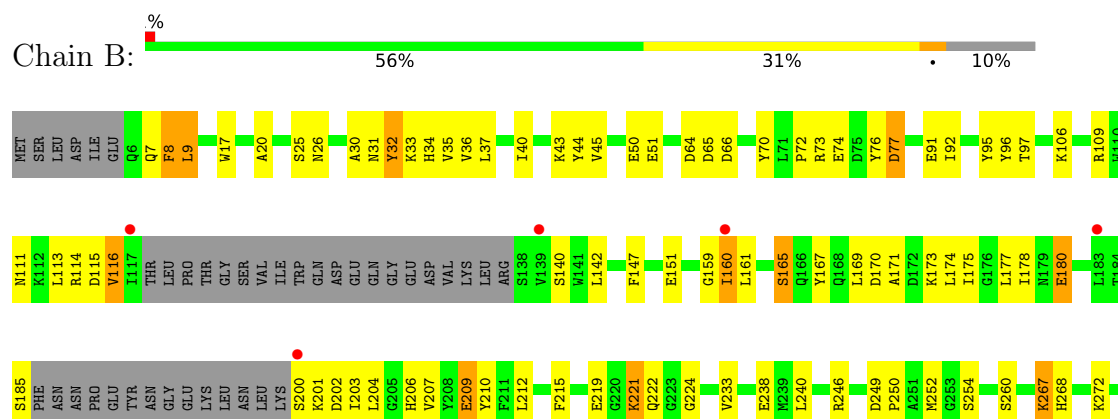
3 Residue-property plots

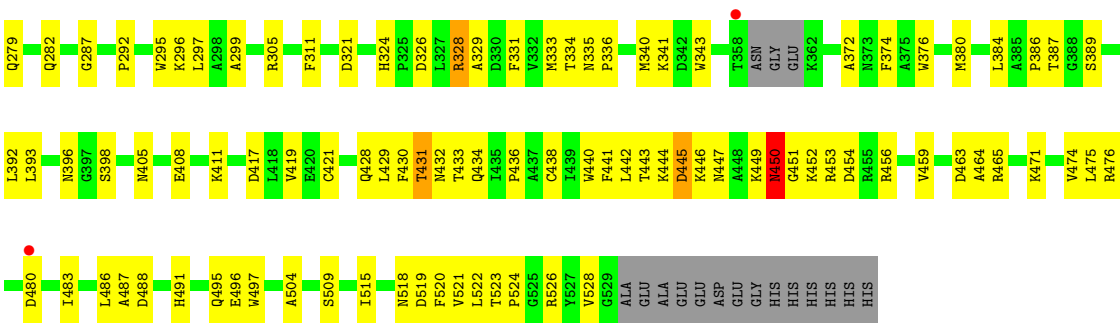
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Type I restriction-modification system methylation subunit



• Molecule 1: Type I restriction-modification system methylation subunit





4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	80.65Å 80.65Å 179.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.55 25.00 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.7 (25.00-2.55) 99.7 (25.00-2.55)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 2.57Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.213 , 0.242 0.209 , 0.234	Depositor DCC
R_{free} test set	1859 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 21.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	0.237 for h,-k,-l	Xtriage
Reported twinning fraction	0.509 for H, K, L 0.491 for K, H, -L	Depositor
Outliers	0 of 37128 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7891	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.28% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.65	0/4027	0.96	5/5453 (0.1%)
1	B	0.65	0/4021	0.95	2/5442 (0.0%)
All	All	0.65	0/8048	0.96	7/10895 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	291	ASN	CA-C-N	5.61	125.07	119.24
1	A	291	ASN	C-N-CA	5.61	125.07	119.24
1	A	437	ALA	N-CA-C	5.37	117.57	109.41
1	A	7	GLN	N-CA-C	5.30	117.58	109.95
1	A	334	THR	N-CA-C	5.07	116.55	108.79
1	B	450	ASN	N-CA-C	5.06	121.58	110.80
1	B	25	SER	N-CA-C	5.02	117.45	109.02

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3934	0	3783	102	0
1	B	3929	0	3764	122	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	5	0	0	1	0
2	B	5	0	0	1	0
3	A	11	0	0	0	0
3	B	7	0	0	0	0
All	All	7891	0	7547	223	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 14.

All (223) 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:334:THR:HG23	1:B:336:PRO:HD3	1.41	1.03
1:B:30:ALA:HA	1:B:219:GLU:OE2	1.61	1.00
1:A:527:TYR:O	1:A:528:VAL:HB	1.56	1.00
1:B:429:LEU:O	1:B:431:THR:HG22	1.78	0.84
1:B:279:GLN:NE2	1:B:282:GLN:HG3	1.94	0.82
1:B:279:GLN:HE21	1:B:282:GLN:HG3	1.44	0.81
1:B:431:THR:HG1	1:B:433:THR:HG1	1.19	0.80
1:B:222:GLN:HB3	1:B:432:ASN:O	1.84	0.77
1:B:160:ILE:HG22	1:B:161:LEU:HD23	1.67	0.76
1:A:139:VAL:HG11	1:A:175:ILE:HD11	1.69	0.74
1:A:431:THR:HG23	1:A:433:THR:H	1.53	0.73
1:B:328:ARG:HD3	2:B:600:SO4:O4	1.88	0.73
1:A:237:VAL:HG22	1:A:333:MET:HE1	1.72	0.71
1:B:454:ASP:OD2	1:B:456:ARG:NH2	2.24	0.71
1:A:210:TYR:CZ	1:A:214:GLN:HG3	2.25	0.70
1:A:410:ARG:HD2	1:A:520:PHE:O	1.91	0.70
1:B:233:VAL:HG12	1:B:260:SER:HB2	1.76	0.67
1:B:36:VAL:O	1:B:40:ILE:HG13	1.94	0.67
1:A:381:LEU:HD21	1:A:418:LEU:HD13	1.77	0.67
1:B:480:ASP:HA	1:B:483:ILE:HB	1.77	0.67
1:A:365:LEU:HD13	1:A:409:ILE:HG12	1.76	0.66
1:B:109:ARG:O	1:B:113:LEU:HD13	1.96	0.65
1:A:157:LEU:HA	1:A:160:ILE:HD12	1.77	0.65
1:B:233:VAL:HG12	1:B:260:SER:CB	2.27	0.65
1:B:171:ALA:O	1:B:175:ILE:HG13	1.97	0.65
1:A:25:SER:OG	1:A:26:ASN:N	2.30	0.64
1:B:160:ILE:HD13	1:B:299:ALA:HB1	1.79	0.63
1:A:527:TYR:O	1:A:528:VAL:CB	2.36	0.63
1:B:449:LYS:O	1:B:451:GLY:N	2.31	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:238:GLU:O	1:B:491:HIS:HE1	1.82	0.62
1:A:232:ILE:HD11	1:A:429:LEU:CD2	2.30	0.62
1:B:453:ARG:HD3	1:B:496:GLU:HA	1.82	0.62
1:B:219:GLU:HG3	1:B:221:LYS:HG2	1.83	0.61
1:B:515:ILE:HG22	1:B:520:PHE:CD1	2.36	0.61
1:B:8:PHE:HD2	1:B:8:PHE:C	2.08	0.61
1:B:238:GLU:O	1:B:491:HIS:CE1	2.54	0.60
1:B:8:PHE:C	1:B:8:PHE:CD2	2.78	0.60
1:A:71:LEU:HD11	1:A:89:GLU:HG3	1.82	0.60
1:A:232:ILE:HD11	1:A:429:LEU:HD21	1.83	0.59
1:B:33:LYS:O	1:B:37:LEU:HG	2.02	0.59
1:B:140:SER:HB3	1:B:165:SER:OG	2.03	0.59
1:B:34:HIS:CE1	1:B:297:LEU:HD21	2.38	0.58
1:B:343:TRP:CD1	1:B:372:ALA:HB2	2.38	0.58
1:B:250:PRO:HD2	1:B:334:THR:HB	1.84	0.58
1:B:174:LEU:O	1:B:178:ILE:HG12	2.04	0.57
1:A:118:THR:HG23	1:A:119:LEU:H	1.69	0.57
1:A:392:LEU:N	1:A:392:LEU:HD12	2.19	0.57
1:A:485:LYS:HE2	1:A:506:PHE:CZ	2.40	0.57
1:A:354:TRP:CD1	1:A:367:PRO:HG3	2.39	0.57
1:A:332:VAL:HG12	1:A:380:MET:HE3	1.87	0.57
1:A:113:LEU:HD22	1:A:142:LEU:HG	1.87	0.56
1:A:232:ILE:HG21	1:A:393:LEU:HD22	1.87	0.56
1:B:384:LEU:HD23	1:B:444:LYS:HG3	1.87	0.56
1:B:465:ARG:O	1:B:476:ARG:NH2	2.38	0.56
1:B:170:ASP:HB3	1:B:173:LYS:HG2	1.87	0.56
1:A:268:HIS:O	1:A:272:LYS:HG2	2.05	0.56
1:A:8:PHE:O	1:A:11:ASP:N	2.39	0.56
1:A:235:LEU:HD11	1:A:239:MET:HE3	1.88	0.55
1:B:240:LEU:HD23	1:B:441:PHE:CZ	2.41	0.55
1:A:328:ARG:NH2	2:A:600:SO4:O1	2.31	0.55
1:B:396:ASN:ND2	1:B:436:PRO:HB2	2.21	0.55
1:A:8:PHE:HD2	1:A:9:LEU:N	2.05	0.54
1:A:346:GLU:HG2	1:B:321:ASP:HA	1.89	0.54
1:A:445:ASP:OD1	1:A:449:LYS:HD2	2.06	0.54
1:A:377:MET:HE1	1:A:440:TRP:CD2	2.42	0.54
1:B:147:PHE:O	1:B:151:GLU:HG3	2.07	0.54
1:A:57:THR:O	1:A:61:GLN:NE2	2.39	0.53
1:A:505:GLY:H	1:A:528:VAL:HG22	1.74	0.53
1:A:213:GLY:O	1:A:216:ALA:HB3	2.08	0.53
1:A:8:PHE:O	1:A:12:LEU:N	2.31	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:LEU:HD13	1:A:440:TRP:HB2	1.91	0.53
1:A:431:THR:HG23	1:A:433:THR:OG1	2.08	0.53
1:B:405:ASN:OD1	1:B:408:GLU:HB2	2.09	0.53
1:A:45:VAL:HG12	1:A:104:VAL:HG22	1.90	0.53
1:B:34:HIS:CE1	1:B:297:LEU:HD11	2.44	0.52
1:B:267:LYS:HE2	1:B:267:LYS:HA	1.90	0.52
1:A:410:ARG:HG3	1:A:440:TRP:HH2	1.74	0.52
1:A:377:MET:HE2	1:A:392:LEU:HD11	1.92	0.52
1:A:514:LEU:O	1:A:518:ASN:ND2	2.34	0.52
1:B:9:LEU:HD12	1:B:206:HIS:CD2	2.45	0.52
1:B:233:VAL:CG1	1:B:260:SER:HB2	2.40	0.52
1:B:44:TYR:C	1:B:44:TYR:CD2	2.88	0.51
1:A:21:ASP:OD2	1:A:24:ARG:HD3	2.10	0.51
1:A:35:VAL:O	1:A:39:LEU:HG	2.11	0.51
1:B:421:CYS:HB3	1:B:441:PHE:HB3	1.92	0.51
1:A:518:ASN:OD1	1:A:526:ARG:HD2	2.11	0.51
1:B:441:PHE:C	1:B:442:LEU:HD12	2.35	0.51
1:A:467:LEU:HD12	1:A:486:LEU:HD11	1.93	0.51
1:A:334:THR:HG23	1:A:336:PRO:HD3	1.93	0.50
1:B:73:ARG:HD3	1:B:73:ARG:C	2.36	0.50
1:B:450:ASN:HD22	1:B:452:LYS:HZ1	1.60	0.50
1:A:112:LYS:O	1:A:116:VAL:HG23	2.11	0.50
1:A:171:ALA:O	1:A:175:ILE:HG12	2.11	0.50
1:B:180:GLU:OE2	1:B:180:GLU:HA	2.10	0.50
1:B:233:VAL:CG1	1:B:260:SER:CB	2.89	0.50
1:B:443:THR:HG22	1:B:445:ASP:H	1.76	0.50
1:A:431:THR:CG2	1:A:433:THR:OG1	2.60	0.49
1:B:246:ARG:HB2	1:B:329:ALA:HA	1.93	0.49
1:A:45:VAL:HG12	1:A:104:VAL:CG2	2.42	0.49
1:A:338:PHE:O	1:A:339:ASN:C	2.55	0.49
1:B:419:VAL:O	1:B:459:VAL:HG22	2.12	0.49
1:A:338:PHE:CE1	1:A:394:LEU:HA	2.48	0.49
1:A:463:ASP:OD1	1:A:465:ARG:HB2	2.12	0.49
1:B:463:ASP:OD2	1:B:465:ARG:NH1	2.46	0.49
1:A:460:LEU:HB2	1:A:493:TRP:CD2	2.47	0.49
1:B:331:PHE:HA	1:B:389:SER:O	2.12	0.49
1:B:92:ILE:HB	1:B:95:TYR:CE2	2.49	0.48
1:A:139:VAL:HG23	1:A:174:LEU:HD22	1.93	0.48
1:A:30:ALA:HA	1:A:33:LYS:HZ2	1.79	0.48
1:A:241:GLU:HG2	1:A:243:TYR:CE2	2.49	0.48
1:A:339:ASN:HA	1:A:371:ASN:OD1	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:453:ARG:HB3	1:B:495:GLN:O	2.14	0.48
1:B:287:GLY:O	1:B:324:HIS:HE1	1.97	0.48
1:A:401:SER:HG	1:A:403:THR:HG1	1.61	0.48
1:A:484:GLN:O	1:A:485:LYS:C	2.57	0.48
1:A:480:ASP:O	1:A:484:GLN:HB2	2.14	0.47
1:B:240:LEU:HD13	1:B:333:MET:HE2	1.96	0.47
1:A:252:MET:HE1	1:A:311:PHE:CD2	2.49	0.47
1:B:34:HIS:ND1	1:B:297:LEU:HD21	2.28	0.47
1:B:209:GLU:CD	1:B:305:ARG:HH21	2.22	0.47
1:A:224:GLY:HA3	1:A:335:ASN:O	2.14	0.47
1:B:20:ALA:O	1:B:32:TYR:OH	2.29	0.47
1:B:70:TYR:CE2	1:B:72:PRO:HA	2.48	0.47
1:B:209:GLU:OE2	1:B:305:ARG:NE	2.44	0.47
1:B:215:PHE:O	1:B:219:GLU:HG2	2.15	0.47
1:A:463:ASP:OD1	1:A:465:ARG:HD3	2.14	0.47
1:A:491:HIS:O	1:A:494:GLN:HB2	2.14	0.47
1:A:391:ALA:HB2	1:A:441:PHE:CD2	2.50	0.47
1:B:518:ASN:O	1:B:519:ASP:HB2	2.14	0.47
1:B:504:ALA:HA	1:B:528:VAL:O	2.14	0.47
1:A:425:LEU:HB3	1:A:429:LEU:HD22	1.96	0.47
1:B:44:TYR:HA	1:B:204:LEU:HD13	1.96	0.47
1:A:463:ASP:OD2	1:A:465:ARG:NH1	2.46	0.46
1:B:111:ASN:O	1:B:114:ARG:HB3	2.15	0.46
1:A:210:TYR:CE2	1:A:214:GLN:HG3	2.50	0.46
1:B:64:ASP:C	1:B:66:ASP:H	2.22	0.45
1:B:92:ILE:HB	1:B:95:TYR:CD2	2.50	0.45
1:B:252:MET:HE1	1:B:311:PHE:CG	2.51	0.45
1:B:252:MET:HE1	1:B:311:PHE:CD2	2.52	0.45
1:A:28:ASP:O	1:A:29:ALA:C	2.58	0.45
1:A:244:LYS:O	1:A:331:PHE:HE2	1.99	0.45
1:A:468:GLY:HA2	1:A:482:ASP:OD2	2.16	0.45
1:B:96:TYR:O	1:B:97:THR:C	2.59	0.45
1:B:222:GLN:HE21	1:B:432:ASN:ND2	2.14	0.45
1:B:30:ALA:CA	1:B:219:GLU:OE2	2.48	0.45
1:A:457:GLY:O	1:A:511:ASP:HA	2.17	0.44
1:B:445:ASP:OD2	1:B:447:ASN:HB2	2.16	0.44
1:A:339:ASN:HA	1:A:371:ASN:CG	2.42	0.44
1:B:471:LYS:HB3	1:B:475:LEU:HB2	1.98	0.44
1:B:487:ALA:O	1:B:488:ASP:C	2.60	0.44
1:A:269:ALA:O	1:A:273:HIS:N	2.49	0.44
1:A:386:PRO:O	1:A:445:ASP:HB2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:ILE:HD13	1:B:207:VAL:HG12	2.00	0.44
1:B:393:LEU:HA	1:B:438:CYS:O	2.18	0.44
1:A:34:HIS:CE1	1:A:293:THR:HG21	2.53	0.44
1:B:376:TRP:O	1:B:380:MET:HG3	2.18	0.43
1:B:389:SER:OG	1:B:443:THR:HG23	2.18	0.43
1:A:235:LEU:CD1	1:A:239:MET:HE3	2.49	0.43
1:B:73:ARG:NH1	1:B:76:TYR:O	2.51	0.43
1:A:411:LYS:HG3	1:A:520:PHE:CE2	2.54	0.43
1:A:112:LYS:O	1:A:112:LYS:HD2	2.18	0.43
1:A:434:GLN:HE21	1:A:434:GLN:HB2	1.60	0.43
1:B:73:ARG:HD3	1:B:73:ARG:O	2.19	0.43
1:B:419:VAL:HA	1:B:442:LEU:HG	2.00	0.43
1:A:327:LEU:O	1:A:383:HIS:ND1	2.51	0.43
1:B:428:GLN:CD	1:B:476:ARG:HG3	2.43	0.43
1:A:392:LEU:HD12	1:A:392:LEU:H	1.83	0.43
1:B:31:ASN:N	1:B:31:ASN:HD22	2.16	0.43
1:B:224:GLY:HA3	1:B:335:ASN:O	2.18	0.43
1:B:374:PHE:CE2	1:B:398:SER:HB2	2.54	0.43
1:A:449:LYS:O	1:A:450:ASN:C	2.62	0.43
1:B:449:LYS:O	1:B:452:LYS:N	2.51	0.43
1:A:516:ARG:HG3	1:A:520:PHE:CE2	2.54	0.43
1:B:26:ASN:O	1:B:26:ASN:CG	2.61	0.43
1:B:173:LYS:O	1:B:177:LEU:HB2	2.19	0.43
1:B:463:ASP:OD1	1:B:463:ASP:C	2.62	0.43
1:A:339:ASN:H	1:A:373:ASN:HD21	1.67	0.43
1:A:425:LEU:O	1:A:426:PRO:C	2.61	0.42
1:B:240:LEU:CD1	1:B:333:MET:HE2	2.49	0.42
1:B:115:ASP:O	1:B:116:VAL:C	2.62	0.42
1:B:464:ALA:HB2	1:B:486:LEU:HD21	2.00	0.42
1:B:474:VAL:HG23	1:B:475:LEU:H	1.85	0.42
1:B:521:VAL:HG21	1:B:526:ARG:HH21	1.85	0.42
1:B:249:ASP:HA	1:B:250:PRO:HD3	1.88	0.42
1:A:243:TYR:OH	1:A:272:LYS:HE3	2.19	0.42
1:A:332:VAL:HG11	1:A:380:MET:HB3	2.01	0.42
1:B:106:LYS:HA	1:B:109:ARG:HD2	2.00	0.42
1:B:386:PRO:HA	1:B:444:LYS:HD3	2.01	0.42
1:B:17:TRP:HB2	1:B:210:TYR:OH	2.19	0.42
1:B:50:GLU:O	1:B:51:GLU:C	2.61	0.42
1:A:162:ASN:ND2	1:A:300:MET:SD	2.92	0.42
1:A:286:TYR:HA	1:A:310:ASN:O	2.19	0.42
1:A:488:ASP:O	1:A:489:THR:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:401:SER:O	1:A:410:ARG:NH2	2.43	0.42
1:B:392:LEU:HD23	1:B:440:TRP:HB2	2.01	0.42
1:B:268:HIS:O	1:B:272:LYS:HB2	2.20	0.41
1:B:335:ASN:CG	1:B:393:LEU:HD23	2.45	0.41
1:B:522:LEU:O	1:B:524:PRO:HD3	2.20	0.41
1:A:207:VAL:O	1:A:207:VAL:CG1	2.67	0.41
1:B:292:PRO:O	1:B:296:LYS:HG3	2.19	0.41
1:B:34:HIS:HE1	1:B:297:LEU:HD11	1.85	0.41
1:B:169:LEU:HG	1:B:173:LYS:HB2	2.02	0.41
1:B:417:ASP:O	1:B:446:LYS:HE3	2.19	0.41
1:A:428:GLN:H	1:A:428:GLN:HG2	1.68	0.41
1:A:80:GLU:H	1:A:80:GLU:CD	2.29	0.41
1:B:449:LYS:HE2	1:B:449:LYS:HB2	1.75	0.41
1:A:414:VAL:O	1:A:415:GLU:C	2.63	0.41
1:B:159:GLY:O	1:B:296:LYS:HE2	2.21	0.41
1:B:200:SER:O	1:B:203:ILE:HB	2.21	0.41
1:B:212:LEU:HD21	1:B:254:SER:HB2	2.03	0.41
1:B:521:VAL:HG21	1:B:526:ARG:NH2	2.35	0.41
1:A:288:GLN:NE2	1:A:316:ALA:O	2.39	0.41
1:A:406:GLU:O	1:A:410:ARG:HB2	2.20	0.41
1:A:348:LEU:O	1:A:351:ASP:HB2	2.21	0.40
1:B:200:SER:HA	1:B:203:ILE:HD12	2.03	0.40
1:A:139:VAL:HG23	1:A:174:LEU:CD2	2.51	0.40
1:A:229:PRO:HB3	1:A:430:PHE:CE2	2.57	0.40
1:B:160:ILE:HG12	1:B:295:TRP:HZ3	1.86	0.40
1:B:200:SER:O	1:B:201:LYS:C	2.64	0.40
1:B:497:TRP:CD1	1:B:509:SER:HB3	2.57	0.40
1:A:17:TRP:CH2	1:A:214:GLN:HB3	2.56	0.40
1:A:240:LEU:HD13	1:A:333:MET:HE3	2.04	0.40
1:B:64:ASP:O	1:B:66:ASP:N	2.55	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	480/544 (88%)	429 (89%)	46 (10%)	5 (1%)	12	17
1	B	479/544 (88%)	414 (86%)	58 (12%)	7 (2%)	8	10
All	All	959/1088 (88%)	843 (88%)	104 (11%)	12 (1%)	9	12

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	528	VAL
1	B	450	ASN
1	A	118	THR
1	B	65	ASP
1	B	74	GLU
1	B	116	VAL
1	B	77	ASP
1	B	430	PHE
1	A	119	LEU
1	A	496	GLU
1	B	32	TYR
1	A	270	ASN

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	413/475 (87%)	372 (90%)	41 (10%)	7	9
1	B	411/475 (86%)	383 (93%)	28 (7%)	14	20
All	All	824/950 (87%)	755 (92%)	69 (8%)	10	14

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLN

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Mol	Chain	Res	Type
1	A	7	GLN
1	A	8	PHE
1	A	12	LEU
1	A	36	VAL
1	A	68	ILE
1	A	79	ASP
1	A	117	ILE
1	A	121	THR
1	A	140	SER
1	A	148	ASP
1	A	149	ASP
1	A	152	LYS
1	A	157	LEU
1	A	166	GLN
1	A	167	TYR
1	A	174	LEU
1	A	177	LEU
1	A	202	ASP
1	A	233	VAL
1	A	240	LEU
1	A	241	GLU
1	A	244	LYS
1	A	267	LYS
1	A	285	VAL
1	A	293	THR
1	A	308	ASP
1	A	334	THR
1	A	340	MET
1	A	342	ASP
1	A	363	ARG
1	A	366	THR
1	A	392	LEU
1	A	411	LYS
1	A	428	GLN
1	A	434	GLN
1	A	443	THR
1	A	449	LYS
1	A	465	ARG
1	A	480	ASP
1	A	481	GLU
1	B	7	GLN
1	B	8	PHE

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Mol	Chain	Res	Type
1	B	9	LEU
1	B	35	VAL
1	B	43	LYS
1	B	45	VAL
1	B	77	ASP
1	B	91	GLU
1	B	142	LEU
1	B	160	ILE
1	B	165	SER
1	B	167	TYR
1	B	180	GLU
1	B	185	SER
1	B	202	ASP
1	B	209	GLU
1	B	221	LYS
1	B	267	LYS
1	B	326	ASP
1	B	328	ARG
1	B	340	MET
1	B	341	LYS
1	B	387	THR
1	B	411	LYS
1	B	431	THR
1	B	434	GLN
1	B	445	ASP
1	B	523	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	34	HIS
1	A	145	ASN
1	A	166	GLN
1	A	214	GLN
1	A	279	GLN
1	A	301	ASN
1	A	396	ASN
1	A	402	ASN
1	A	405	ASN
1	A	416	GLN
1	A	432	ASN

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Mol	Chain	Res	Type
1	A	434	GLN
1	A	450	ASN
1	A	458	GLN
1	A	503	GLN
1	B	26	ASN
1	B	31	ASN
1	B	34	HIS
1	B	83	GLN
1	B	145	ASN
1	B	206	HIS
1	B	279	GLN
1	B	301	ASN
1	B	315	ASN
1	B	324	HIS
1	B	404	ASN
1	B	432	ASN
1	B	450	ASN
1	B	458	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
2	SO4	A	600	-	4,4,4	0.57	0	6,6,6	0.46	0
2	SO4	B	600	-	4,4,4	0.64	0	6,6,6	0.54	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	SO4	1	0
2	B	600	SO4	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	488/544 (89%)	0.11	9 (1%) 67 68	30, 52, 68, 80	0
1	B	487/544 (89%)	0.15	7 (1%) 73 74	27, 51, 70, 82	0
All	All	975/1088 (89%)	0.13	16 (1%) 70 71	27, 52, 68, 82	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	433	THR	2.9
1	A	220	GLY	2.8
1	A	218	ALA	2.8
1	B	117	ILE	2.7
1	A	480	ASP	2.5
1	B	183	LEU	2.4
1	A	225	GLN	2.3
1	A	182	SER	2.2
1	B	358	THR	2.2
1	B	160	ILE	2.2
1	B	200	SER	2.2
1	A	441	PHE	2.1
1	B	480	ASP	2.1
1	A	117	ILE	2.0
1	B	139	VAL	2.0
1	A	529	GLY	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 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	SO4	A	600	5/5	0.89	0.31	44,44,48,48	0
2	SO4	B	600	5/5	0.91	0.31	39,40,44,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.