



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:18 PM UTC

PDB ID : 4IA5 / pdb_00004ia5
Title : Hydratase from Lactobacillus acidophilus - SeMet derivative (apo LAH)
Authors : Khoshnevis, S.; Neumann, P.; Ficner, R.
Deposited on : 2012-12-06
Resolution : 2.22 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

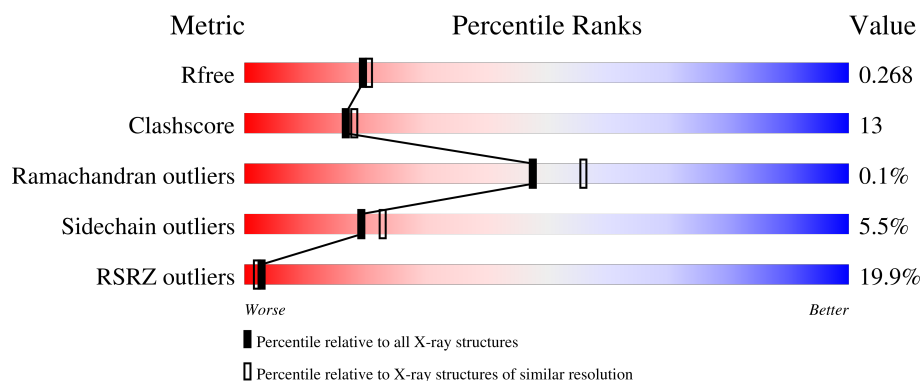
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	591	9% 76% 20% • •
1	B	591	29% 65% 28% 5% •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	GOL	B	608	-	-	X	-

2 Entry composition [i](#)

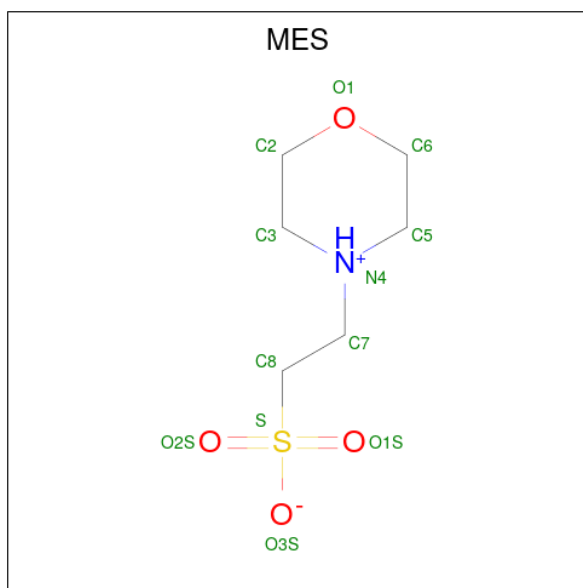
There are 8 unique types of molecules in this entry. The entry contains 9939 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Myosin-crossreactive antigen.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	577	Total	C	N	O	S	Se	0	0	0
			4679	2992	771	892	6	18			
1	B	579	Total	C	N	O	S	Se	0	1	0
			4696	3002	774	895	6	19			

- Molecule 2 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total K 2 2	0	0

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 8 6 2	0	0
4	A	1	Total C O 8 6 2	0	0
4	A	1	Total C O 8 6 2	0	0
4	B	1	Total C O 8 6 2	0	0
4	B	1	Total C O 8 6 2	0	0

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Na 1 1	0	0

- Molecule 6 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

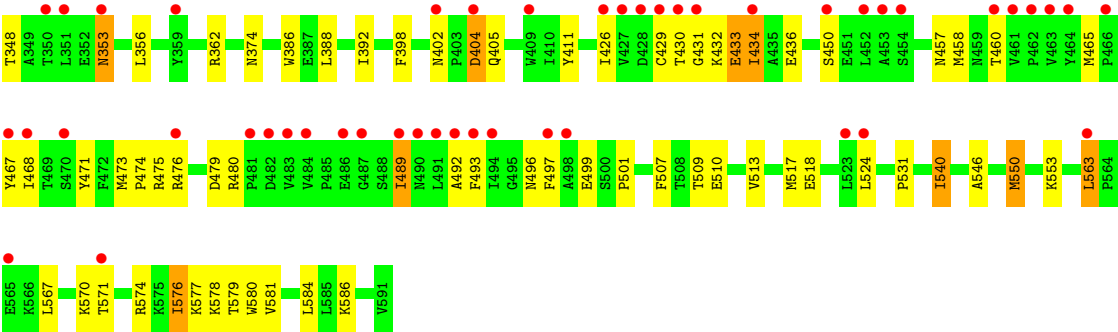
- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	209	Total 209	O 209	0	0
8	B	264	Total 265	O 265	0	1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	168.75Å 78.97Å 108.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.79 – 2.22 44.79 – 2.22	Depositor EDS
% Data completeness (in resolution range)	88.4 (44.79-2.22) 92.9 (44.79-2.22)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.86 (at 2.22Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.214 , 0.259 0.233 , 0.268	Depositor DCC
R_{free} test set	3370 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.615	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 59.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9939	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% 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: MES, MPD, PO4, K, GOL, NA

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	0.43	0/4781	0.84	1/6448 (0.0%)
1	B	0.53	2/4798 (0.0%)	0.95	15/6470 (0.2%)
All	All	0.48	2/9579 (0.0%)	0.89	16/12918 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	92	ASP	C-O	5.47	1.30	1.24
1	B	305	ALA	CA-C	5.07	1.57	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	489	ILE	N-CA-C	8.45	119.22	112.12
1	B	97	ILE	N-CA-C	7.22	116.56	108.05
1	B	303	THR	CA-C-N	7.02	128.62	119.84
1	B	303	THR	C-N-CA	7.02	128.62	119.84
1	B	330	HIS	CA-C-N	7.00	126.98	119.28
1	B	330	HIS	C-N-CA	7.00	126.98	119.28
1	B	268	MSE	N-CA-C	6.17	116.71	108.38
1	B	277	ILE	N-CA-C	5.38	116.49	107.18
1	B	305	ALA	CA-C-N	5.37	127.50	120.25
1	B	305	ALA	C-N-CA	5.37	127.50	120.25
1	B	279	LEU	N-CA-C	5.37	117.83	108.76
1	B	306	PRO	N-CA-C	5.33	122.19	111.69
1	B	86	HIS	N-CA-C	5.28	118.56	112.97
1	A	324	GLN	N-CA-C	5.17	116.99	111.36
1	B	331	PRO	N-CA-C	5.14	120.59	113.65
1	B	251	ASN	N-CA-C	5.04	115.80	108.14

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	4679	0	4543	90	0
1	B	4696	0	4558	151	0
2	A	12	0	12	0	0
2	B	12	0	12	0	0
3	A	2	0	0	0	0
4	A	24	0	42	5	0
4	B	16	0	28	6	0
5	B	1	0	0	0	0
6	B	5	0	0	0	0
7	B	18	0	24	5	0
8	A	209	0	0	4	0
8	B	265	0	0	11	0
All	All	9939	0	9219	240	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 13.

All (240) 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:52:ILE:HB	1:A:241:VAL:HG23	1.57	0.87
1:B:47[A]:MSE:HE3	1:B:524:LEU:HD11	1.63	0.81
1:A:186:PHE:HE1	4:A:604:MPD:H31	1.46	0.81
1:B:28:ILE:HD11	1:B:287:VAL:HG22	1.63	0.80
1:B:266:ILE:HB	1:B:277:ILE:HG13	1.65	0.78
1:B:429:CYS:HB3	1:B:433:GLU:HB2	1.64	0.77
1:B:36:SER:HB3	1:B:93:MSE:HE1	1.69	0.75
1:A:23:LYS:O	1:A:51:ASN:ND2	2.20	0.73
1:A:158:LYS:NZ	1:A:477:ASP:OD2	2.19	0.72
1:B:92:ASP:O	1:B:95:ARG:HG3	1.90	0.71
1:B:196:GLU:OE1	1:B:475:ARG:NH2	2.24	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:LEU:HB3	1:A:241:VAL:HG11	1.73	0.71
1:B:49:GLY:HA2	1:B:52:ILE:HD12	1.73	0.71
1:B:249:VAL:HA	1:B:268:MSE:HB2	1.70	0.71
1:B:41:LEU:HB3	1:B:47[B]:MSE:HE2	1.73	0.70
1:B:290:GLY:H	1:B:314:SER:HB3	1.56	0.70
1:B:338:LEU:HG	1:B:473:MSE:HE3	1.74	0.70
1:B:47[B]:MSE:HE3	1:B:52:ILE:HG12	1.73	0.69
1:B:252:VAL:HG22	1:B:266:ILE:HG12	1.73	0.69
1:B:411:TYR:HB2	4:B:604:MPD:H11	1.73	0.69
1:B:266:ILE:HD11	1:B:285:VAL:HG21	1.75	0.68
1:B:236:LEU:HD23	1:B:241:VAL:HG21	1.75	0.68
1:B:479:ASP:O	8:B:961:HOH:O	2.14	0.65
1:A:567:LEU:HD12	1:A:570:LYS:HZ2	1.62	0.64
1:A:26:TYR:HB2	1:A:285:VAL:HG22	1.79	0.64
1:A:359:TYR:O	1:A:362:ARG:HG2	1.97	0.64
1:B:208:ILE:HG21	1:B:546:ALA:HA	1.80	0.64
1:B:251:ASN:OD1	1:B:252:VAL:N	2.31	0.63
1:B:307:ILE:O	7:B:608:GOL:H31	1.99	0.62
1:A:538:TYR:HD1	1:B:540:ILE:HD11	1.63	0.61
1:B:540:ILE:HD13	1:B:584:LEU:HD13	1.82	0.61
1:B:228:MSE:HE1	1:B:509:THR:HG21	1.82	0.61
1:B:26:TYR:HB2	1:B:285:VAL:HG22	1.81	0.61
1:B:313:ASP:O	1:B:317:LEU:HG	2.01	0.61
1:A:254:VAL:H	1:A:324:GLN:HE22	1.49	0.61
1:B:52:ILE:HB	1:B:241:VAL:HG13	1.82	0.61
1:A:82:GLU:HG3	1:A:217:LEU:HD22	1.82	0.60
1:A:82:GLU:OE2	4:A:604:MPD:O2	2.18	0.60
1:B:41:LEU:HD13	1:B:47[B]:MSE:HE1	1.83	0.60
1:A:366:ARG:NH1	1:A:374:ASN:O	2.31	0.60
1:B:303:THR:N	1:B:430:THR:OG1	2.35	0.59
1:A:154:MSE:HE3	1:B:579:THR:HG21	1.84	0.59
1:A:566:LYS:O	1:A:570:LYS:HG2	2.03	0.59
1:A:340:GLU:OE2	1:A:425:ARG:NH1	2.36	0.59
1:A:561:MSE:HE3	1:A:566:LYS:HD3	1.85	0.58
1:A:252:VAL:HG13	1:A:321:LEU:HD21	1.86	0.58
1:B:36:SER:CB	1:B:93:MSE:HE1	2.32	0.58
1:B:185:MSE:HE3	1:B:214:PHE:HE1	1.68	0.58
1:B:73:ALA:HB1	1:B:398:PHE:CD2	2.39	0.57
1:A:297:THR:HB	1:A:306:PRO:HD2	1.86	0.57
1:A:262:ILE:HD13	1:A:281:HIS:CE1	2.40	0.57
1:B:54:ILE:HD11	1:B:243:PHE:CD2	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:TYR:OH	1:B:518:GLU:OE2	2.18	0.56
1:B:32:LEU:N	8:B:834:HOH:O	2.37	0.56
1:A:186:PHE:CE1	4:A:604:MPD:H31	2.35	0.56
1:B:576:ILE:HD13	1:B:581:VAL:HG12	1.87	0.56
1:A:254:VAL:HG22	1:A:263:ALA:HA	1.88	0.56
1:B:289:ASN:ND2	8:B:896:HOH:O	2.36	0.55
1:B:332:ASP:OD1	1:B:332:ASP:N	2.37	0.55
1:A:205:ILE:O	1:A:208:ILE:HG12	2.06	0.55
1:B:26:TYR:CE2	1:B:283:ASP:HB3	2.41	0.55
1:A:80:GLY:HA3	1:A:224:GLN:HB2	1.88	0.55
1:A:11:PHE:CG	1:B:95:ARG:HG2	2.42	0.54
1:B:430:THR:H	1:B:433:GLU:HG3	1.71	0.54
1:B:432:LYS:HE3	1:B:436:GLU:OE2	2.07	0.54
1:B:286:PHE:CD1	1:B:492:ALA:HB3	2.42	0.54
1:B:208:ILE:HD11	1:B:550:MSE:HG3	1.88	0.54
1:B:232:LEU:O	1:B:236:LEU:HD12	2.08	0.54
1:B:318:TRP:CD2	1:B:331:PRO:HB3	2.43	0.53
1:A:330:HIS:O	1:A:333:VAL:HG12	2.09	0.53
1:A:87:PHE:CG	1:A:510:GLU:HB2	2.42	0.53
1:A:292:ILE:HD13	1:A:471:TYR:CE2	2.43	0.53
1:B:82:GLU:HG3	1:B:217:LEU:HD22	1.91	0.53
1:A:267:VAL:HG12	1:A:276:GLU:HG2	1.91	0.53
1:A:416:ASP:HA	1:A:425:ARG:HD2	1.89	0.53
1:B:74:GLY:HA2	1:B:348:THR:HG21	1.90	0.53
1:B:57:GLU:HA	1:B:247:CYS:O	2.09	0.52
1:B:402:ASN:ND2	1:B:404:ASP:OD1	2.42	0.52
1:B:252:VAL:O	1:B:324:GLN:NE2	2.42	0.52
1:B:87:PHE:CG	1:B:510:GLU:HB2	2.45	0.52
1:B:166:GLU:HB2	1:B:190:LYS:O	2.10	0.52
1:A:246:ASP:OD1	1:A:248:HIS:NE2	2.34	0.52
1:B:22:LYS:O	1:B:282:ASN:HB3	2.10	0.52
1:B:311:LYS:NZ	1:B:336:GLU:OE1	2.21	0.52
1:B:374:ASN:HD22	1:B:392:ILE:HG22	1.73	0.52
1:B:80:GLY:HA3	1:B:224:GLN:HB2	1.90	0.52
1:A:87:PHE:CD2	1:A:510:GLU:HB2	2.45	0.52
1:A:264:LYS:O	1:A:279:LEU:N	2.40	0.52
1:B:27:ILE:HB	1:B:54:ILE:HG22	1.91	0.52
1:B:262:ILE:HG12	1:B:489:ILE:HD12	1.92	0.51
1:B:204:PHE:CZ	4:B:605:MPD:HM2	2.46	0.51
1:B:579:THR:HG22	1:B:580:TRP:N	2.26	0.51
1:B:93:MSE:HE3	1:B:513:VAL:HG13	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:184:THR:O	4:B:604:MPD:O4	2.28	0.51
1:B:307:ILE:HB	7:B:608:GOL:H31	1.93	0.50
1:B:274:ASP:OD1	1:B:275:LYS:N	2.44	0.50
1:A:544:LEU:HA	1:A:547:MSE:HE2	1.94	0.50
1:A:76:VAL:HA	1:A:409:TRP:CH2	2.47	0.50
1:A:82:GLU:H	1:A:82:GLU:CD	2.20	0.50
1:A:50:GLU:HG2	8:A:897:HOH:O	2.10	0.50
1:A:266:ILE:HD12	1:A:277:ILE:HD11	1.94	0.50
1:A:133:ARG:HE	1:A:137:ASP:HB3	1.77	0.50
1:B:30:SER:OG	1:B:56:GLU:HG2	2.12	0.50
1:B:310:ALA:O	1:B:467:TYR:OH	2.16	0.49
1:B:37:THR:HG21	1:B:286:PHE:CD2	2.47	0.49
1:B:81:ARG:NH2	8:B:804:HOH:O	2.30	0.49
1:B:465:MSE:SE	1:B:468:ILE:HD13	2.62	0.49
1:B:38:ALA:HA	1:B:41:LEU:HD12	1.95	0.49
1:B:321:LEU:HB3	1:B:328:PHE:CD2	2.47	0.49
1:A:270:GLN:O	1:A:273:LYS:HG3	2.12	0.48
1:A:279:LEU:HD22	1:A:283:ASP:CG	2.38	0.48
1:B:307:ILE:HB	7:B:608:GOL:C3	2.43	0.48
1:A:161:GLU:HG2	1:A:476:ARG:HG2	1.94	0.48
1:B:480:ARG:NE	1:B:499:GLU:OE1	2.43	0.48
1:A:471:TYR:CZ	1:A:496:ASN:HB2	2.48	0.48
1:B:292:ILE:HD11	7:B:606:GOL:H12	1.95	0.48
1:A:26:TYR:CD1	1:A:53:HIS:HB2	2.49	0.48
1:A:147:ASN:ND2	8:A:829:HOH:O	2.47	0.48
1:B:23:LYS:HA	1:B:282:ASN:HB3	1.94	0.48
1:B:251:ASN:HB3	1:B:267:VAL:CG2	2.44	0.48
1:B:426:ILE:HG13	1:B:434:ILE:HD11	1.95	0.47
1:A:234:ALA:O	1:A:238:ASP:HB2	2.13	0.47
1:A:368:LEU:HA	1:A:374:ASN:HD21	1.80	0.47
1:A:423:LYS:HA	1:A:423:LYS:HD3	1.68	0.47
1:B:3:TYR:OH	1:B:501:PRO:HG2	2.14	0.47
1:B:26:TYR:HE1	1:B:53:HIS:HD1	1.62	0.47
1:B:193:SER:HB2	1:B:474:PRO:HA	1.96	0.47
1:A:51:ASN:HA	8:A:898:HOH:O	2.13	0.47
1:A:421:TYR:CE1	1:A:446:GLU:HG2	2.49	0.47
1:B:320:ASN:O	1:B:324:GLN:HG3	2.15	0.47
1:B:28:ILE:HG13	1:B:287:VAL:HA	1.96	0.47
1:B:432:LYS:HD2	8:B:924:HOH:O	2.14	0.47
1:A:24:SER:HB3	1:A:51:ASN:OD1	2.14	0.47
1:B:75:PHE:HB3	1:B:78:ARG:HG3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:421:TYR:HE1	1:A:446:GLU:HG2	1.80	0.46
1:B:280:THR:HG22	1:B:282:ASN:H	1.80	0.46
1:B:476:ARG:O	1:B:479:ASP:HB2	2.15	0.46
1:B:292:ILE:HD13	1:B:471:TYR:CE2	2.51	0.46
1:B:330:HIS:HB2	1:B:479:ASP:OD1	2.16	0.46
1:B:23:LYS:HG2	1:B:282:ASN:HA	1.97	0.46
1:B:307:ILE:H	7:B:608:GOL:H12	1.80	0.46
1:B:386:TRP:HE3	1:B:388:LEU:HD13	1.81	0.46
1:B:546:ALA:O	1:B:550:MSE:HB2	2.15	0.46
1:B:233:LEU:HD22	1:B:243:PHE:CE2	2.51	0.46
1:B:431:GLY:HA3	1:B:460:THR:O	2.15	0.46
1:B:174:PHE:CD1	1:B:190:LYS:HE2	2.51	0.46
1:B:356:LEU:HD13	1:B:458:MSE:SE	2.66	0.46
1:A:130:ARG:HD3	1:A:440:TYR:CZ	2.51	0.45
1:A:35:LEU:O	1:A:39:VAL:HG23	2.16	0.45
1:B:430:THR:HG22	1:B:433:GLU:CG	2.46	0.45
1:A:24:SER:OG	1:A:283:ASP:OD1	2.34	0.45
1:B:318:TRP:CG	1:B:331:PRO:HB3	2.51	0.45
1:B:318:TRP:HZ2	8:B:961:HOH:O	1.99	0.45
1:B:335:CYS:HB2	1:B:467:TYR:CE2	2.51	0.45
1:B:567:LEU:O	1:B:571:THR:HG23	2.16	0.45
1:B:33:ALA:HB3	1:B:288:THR:HG23	1.98	0.45
1:B:430:THR:HG22	1:B:433:GLU:HG3	1.99	0.45
1:A:417:THR:O	1:A:425:ARG:HA	2.16	0.45
1:A:353:ASN:OD1	1:A:355:LYS:HG3	2.16	0.45
1:A:485:PRO:HG2	1:A:488:SER:HB2	1.99	0.45
1:B:47[B]:MSE:HE2	1:B:47[B]:MSE:HB3	1.82	0.45
1:B:290:GLY:N	1:B:314:SER:HB3	2.26	0.45
1:A:24:SER:HA	1:A:51:ASN:ND2	2.31	0.45
1:A:204:PHE:CZ	4:A:604:MPD:HM2	2.52	0.45
1:A:251:ASN:OD1	1:A:252:VAL:N	2.49	0.45
1:B:59:PRO:HB3	1:B:246:ASP:HB2	1.99	0.45
1:A:82:GLU:HB3	1:A:219:PHE:CE2	2.52	0.44
1:B:91:TRP:HA	1:B:91:TRP:CE3	2.51	0.44
1:B:226:GLU:HG3	8:B:956:HOH:O	2.18	0.44
1:B:471:TYR:CZ	1:B:496:ASN:HB2	2.53	0.44
1:B:430:THR:N	1:B:433:GLU:HG3	2.32	0.44
1:A:567:LEU:HD12	1:A:570:LYS:NZ	2.32	0.44
1:B:47[B]:MSE:HE3	1:B:52:ILE:CG1	2.43	0.44
1:B:144:LYS:HG2	8:B:864:HOH:O	2.17	0.44
1:A:578:LYS:HD2	1:A:578:LYS:HA	1.73	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:220:ASN:HD21	1:B:228:MSE:SE	2.51	0.43
1:B:553:LYS:NZ	8:B:937:HOH:O	2.51	0.43
1:A:26:TYR:HD1	1:A:53:HIS:HB2	1.83	0.43
1:A:278:ASP:O	1:A:279:LEU:HD23	2.18	0.43
1:B:15:LYS:HE3	1:B:15:LYS:HB2	1.86	0.43
1:B:286:PHE:CE1	1:B:492:ALA:HB3	2.54	0.43
1:A:48:LYS:HZ3	1:A:50:GLU:CD	2.25	0.43
1:B:28:ILE:CD1	1:B:287:VAL:HG22	2.41	0.43
1:B:237:LYS:HD3	1:B:243:PHE:CE1	2.53	0.43
1:B:450:SER:HA	8:B:957:HOH:O	2.19	0.43
1:B:23:LYS:HB3	1:B:47[A]:MSE:HE2	1.99	0.43
1:B:54:ILE:HD11	1:B:243:PHE:CE2	2.54	0.43
1:B:254:VAL:HG22	1:B:263:ALA:HA	2.00	0.43
1:A:76:VAL:HB	1:A:463:VAL:HG21	2.00	0.42
1:A:236:LEU:HB3	1:A:241:VAL:CG1	2.46	0.42
1:B:24:SER:N	1:B:282:ASN:O	2.38	0.42
1:B:318:TRP:CH2	1:B:497:PHE:HB3	2.53	0.42
1:A:236:LEU:O	1:A:241:VAL:HG12	2.19	0.42
1:A:340:GLU:HG3	1:A:416:ASP:HB3	2.00	0.42
1:A:368:LEU:HA	1:A:374:ASN:ND2	2.33	0.42
1:B:339:PRO:HD2	1:B:342:SER:HG	1.84	0.42
1:B:97:ILE:HA	1:B:98:PRO:HD3	1.94	0.42
1:B:517:MSE:SE	1:B:531:PRO:HD2	2.70	0.42
1:B:255:ASP:OD2	1:B:264:LYS:HD3	2.20	0.42
1:B:321:LEU:HD13	1:B:493:PHE:CZ	2.54	0.42
1:A:160:ILE:HB	1:A:194:LEU:HD23	2.02	0.42
1:B:244:GLU:HB3	1:B:247:CYS:SG	2.60	0.42
1:B:563:LEU:H	1:B:563:LEU:HD12	1.84	0.42
1:A:236:LEU:C	1:A:241:VAL:HG12	2.45	0.42
1:A:237:LYS:HE3	1:A:237:LYS:HB2	1.89	0.42
1:A:270:GLN:HG2	1:A:271:ASN:OD1	2.20	0.42
1:A:432:LYS:HE3	1:A:454:SER:HA	2.02	0.42
1:A:356:LEU:HD22	1:A:439:LEU:HD22	2.02	0.41
1:A:523:LEU:HD23	1:A:523:LEU:HA	1.90	0.41
1:A:540:ILE:O	1:A:544:LEU:HG	2.21	0.41
1:B:75:PHE:HD1	1:B:77:VAL:HG12	1.85	0.41
1:B:169:PHE:CZ	1:B:197:MSE:HE2	2.55	0.41
1:B:274:ASP:C	1:B:275:LYS:HG2	2.44	0.41
1:A:202:MSE:HE3	1:A:537:ILE:HG12	2.01	0.41
1:A:576:ILE:O	1:A:579:THR:OG1	2.27	0.41
1:A:97:ILE:HA	1:A:98:PRO:HD3	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:GLU:H	1:B:82:GLU:CD	2.27	0.41
1:B:402:ASN:H	1:B:405:GLN:NE2	2.18	0.41
1:A:107:TYR:OH	1:A:227:SER:HB3	2.20	0.41
1:A:253:GLU:HA	1:A:324:GLN:HE22	1.85	0.41
1:A:397:HIS:HD2	8:A:742:HOH:O	2.02	0.41
1:A:87:PHE:HB3	1:A:90:LEU:HB3	2.02	0.41
1:A:25:ALA:HA	1:A:284:ILE:O	2.20	0.41
1:A:471:TYR:CE1	1:A:496:ASN:HB2	2.56	0.41
1:B:191:TRP:CZ3	1:B:341:ARG:HG2	2.56	0.41
1:B:321:LEU:HD13	1:B:493:PHE:HZ	1.85	0.41
1:B:353:ASN:OD1	1:B:457:ASN:HA	2.20	0.41
4:B:605:MPD:HM1	4:B:605:MPD:H4	1.90	0.41
1:B:84:GLU:HG2	4:B:605:MPD:H32	2.03	0.41
1:B:277:ILE:HG12	8:B:962:HOH:O	2.19	0.41
1:B:411:TYR:HB2	4:B:604:MPD:C1	2.46	0.41
1:B:251:ASN:CG	1:B:252:VAL:N	2.79	0.40
1:A:58:LEU:HA	1:A:59:PRO:HD3	1.92	0.40
1:B:4:SER:O	1:B:531:PRO:HA	2.21	0.40
1:B:107:TYR:OH	1:B:227:SER:HB3	2.21	0.40
4:A:606:MPD:H4	4:A:606:MPD:HM1	1.78	0.40
1:B:24:SER:OG	1:B:51:ASN:O	2.32	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	573/591 (97%)	563 (98%)	9 (2%)	1 (0%)	43	50
1	B	576/591 (98%)	565 (98%)	11 (2%)	0	100	100
All	All	1149/1182 (97%)	1128 (98%)	20 (2%)	1 (0%)	48	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	423	LYS

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	511/500 (102%)	489 (96%)	22 (4%)	26	33
1	B	512/500 (102%)	477 (93%)	35 (7%)	14	16
All	All	1023/1000 (102%)	966 (94%)	57 (6%)	19	22

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	24	SER
1	A	52	ILE
1	A	75	PHE
1	A	128	TYR
1	A	159	GLU
1	A	257	GLU
1	A	269	THR
1	A	271	ASN
1	A	273	LYS
1	A	278	ASP
1	A	323	LYS
1	A	332	ASP
1	A	382	VAL
1	A	447	SER
1	A	448	GLN
1	A	454	SER
1	A	563	LEU
1	A	570	LYS
1	A	571	THR
1	A	578	LYS
1	A	587	GLU

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Mol	Chain	Res	Type
1	B	1	MSE
1	B	28	ILE
1	B	30	SER
1	B	47[A]	MSE
1	B	47[B]	MSE
1	B	48	LYS
1	B	50	GLU
1	B	56	GLU
1	B	84	GLU
1	B	95	ARG
1	B	151	LYS
1	B	197	MSE
1	B	226	GLU
1	B	247	CYS
1	B	273	LYS
1	B	277	ILE
1	B	288	THR
1	B	293	THR
1	B	314	SER
1	B	332	ASP
1	B	353	ASN
1	B	362	ARG
1	B	404	ASP
1	B	433	GLU
1	B	434	ILE
1	B	507	PHE
1	B	540	ILE
1	B	550	MSE
1	B	563	LEU
1	B	570	LYS
1	B	574	ARG
1	B	576	ILE
1	B	577	LYS
1	B	578	LYS
1	B	586	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	ASN
1	A	256	HIS
1	A	324	GLN

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Mol	Chain	Res	Type
1	A	374	ASN
1	A	393	HIS
1	A	459	ASN
1	B	139	GLN
1	B	147	ASN
1	B	223	ASN
1	B	385	ASN
1	B	393	HIS
1	B	496	ASN

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](#)

Of 14 ligands modelled in this entry, 3 are monoatomic - leaving 11 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	$\# Z > 2$
4	MPD	A	605	-	7,7,7	0.31	0	9,10,10	0.23	0
4	MPD	B	605	-	7,7,7	0.26	0	9,10,10	0.50	0
6	PO4	B	603	-	4,4,4	1.04	0	6,6,6	0.52	0
7	GOL	B	606	-	5,5,5	0.34	0	5,5,5	0.33	0
7	GOL	B	608	-	5,5,5	0.38	0	5,5,5	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MPD	A	606	-	7,7,7	0.28	0	9,10,10	0.27	0
7	GOL	B	607	-	5,5,5	0.35	0	5,5,5	0.36	0
2	MES	B	602	-	12,12,12	2.17	1 (8%)	15,16,16	2.45	7 (46%)
2	MES	A	601	-	12,12,12	2.22	1 (8%)	15,16,16	1.96	6 (40%)
4	MPD	B	604	-	7,7,7	0.21	0	9,10,10	0.44	0
4	MPD	A	604	-	7,7,7	0.32	0	9,10,10	0.54	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
4	MPD	A	605	-	-	3/5/5/5	-
4	MPD	B	605	-	-	2/5/5/5	-
7	GOL	B	606	-	-	0/4/4/4	-
7	GOL	B	608	-	-	1/4/4/4	-
4	MPD	A	606	-	-	2/5/5/5	-
7	GOL	B	607	-	-	3/4/4/4	-
2	MES	B	602	-	-	1/6/14/14	0/1/1/1
2	MES	A	601	-	-	4/6/14/14	0/1/1/1
4	MPD	B	604	-	-	1/5/5/5	-
4	MPD	A	604	-	-	2/5/5/5	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	MES	C8-S	-7.31	1.67	1.77
2	B	602	MES	C8-S	-7.17	1.67	1.77

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	602	MES	C5-N4-C3	5.43	120.53	108.84
2	A	601	MES	C5-N4-C3	4.23	117.95	108.84
2	B	602	MES	O1S-S-C8	3.99	112.76	106.73
2	B	602	MES	C7-N4-C3	3.35	120.17	111.24
2	A	601	MES	O2S-S-C8	3.14	111.47	106.73
2	B	602	MES	C7-N4-C5	2.97	119.16	111.24
2	A	601	MES	C6-C5-N4	-2.43	106.43	110.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	MES	C7-N4-C5	2.41	117.67	111.24
2	B	602	MES	C6-C5-N4	-2.33	106.59	110.12
2	B	602	MES	O2S-S-C8	2.30	110.20	106.73
2	A	601	MES	O1S-S-C8	2.27	110.16	106.73
2	A	601	MES	C7-N4-C3	2.15	116.96	111.24
2	B	602	MES	C2-C3-N4	-2.14	106.87	110.12

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	MES	C8-C7-N4-C3
4	A	604	MPD	C2-C3-C4-C5
4	B	605	MPD	C2-C3-C4-O4
4	B	605	MPD	C2-C3-C4-C5
7	B	607	GOL	C1-C2-C3-O3
2	B	602	MES	C8-C7-N4-C5
2	A	601	MES	C7-C8-S-O3S
7	B	607	GOL	O2-C2-C3-O3
7	B	608	GOL	O1-C1-C2-O2
2	A	601	MES	C7-C8-S-O1S
2	A	601	MES	C7-C8-S-O2S
4	A	604	MPD	O2-C2-C3-C4
4	A	606	MPD	C2-C3-C4-O4
4	A	605	MPD	C1-C2-C3-C4
4	A	605	MPD	CM-C2-C3-C4
4	A	606	MPD	C2-C3-C4-C5
4	A	605	MPD	O2-C2-C3-C4
4	B	604	MPD	C1-C2-C3-C4
7	B	607	GOL	O1-C1-C2-C3

There are no ring outliers.

6 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	605	MPD	3	0
7	B	606	GOL	1	0
7	B	608	GOL	4	0
4	A	606	MPD	1	0
4	B	604	MPD	3	0
4	A	604	MPD	4	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	559/591 (94%)	0.84	53 (9%) 14 11	29, 61, 99, 127	0
1	B	561/591 (94%)	1.32	170 (30%) 1 1	25, 67, 109, 127	4 (0%)
All	All	1120/1182 (94%)	1.08	223 (19%) 3 2	25, 63, 105, 127	4 (0%)

All (223) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	267	VAL	5.9
1	B	317	LEU	5.5
1	B	305	ALA	5.4
1	B	327	ALA	5.0
1	B	60	VAL	5.0
1	B	279	LEU	4.7
1	B	307	ILE	4.7
1	B	286	PHE	4.6
1	B	266	ILE	4.5
1	B	309	ASN	4.5
1	B	254	VAL	4.4
1	B	314	SER	4.4
1	B	427	VAL	4.4
1	A	76	VAL	4.3
1	B	245	TYR	4.3
1	B	249	VAL	4.3
1	B	333	VAL	4.3
1	B	298	TYR	4.3
1	B	296	SER	4.2
1	B	304	PRO	4.2
1	B	31	GLY	4.1
1	B	277	ILE	4.1
1	A	60	VAL	4.1
1	B	310	ALA	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	78	ARG	4.1
1	B	75	PHE	4.1
1	A	75	PHE	4.1
1	A	489	ILE	4.1
1	B	73	ALA	4.0
1	B	288	THR	4.0
1	B	54	ILE	4.0
1	A	556	LEU	4.0
1	B	321	LEU	4.0
1	B	74	GLY	3.9
1	B	467	TYR	3.9
1	B	262	ILE	3.8
1	B	338	LEU	3.8
1	B	301	GLN	3.8
1	A	277	ILE	3.8
1	B	489	ILE	3.8
1	B	303	THR	3.8
1	B	27	ILE	3.8
1	A	572	GLY	3.8
1	B	318	TRP	3.7
1	B	491	LEU	3.7
1	B	252	VAL	3.7
1	A	281	HIS	3.7
1	B	306	PRO	3.7
1	B	497	PHE	3.6
1	B	271	ASN	3.6
1	B	272	GLY	3.6
1	B	493	PHE	3.6
1	A	333	VAL	3.6
1	A	262	ILE	3.5
1	B	292	ILE	3.5
1	B	329	GLY	3.5
1	A	258	GLY	3.5
1	B	315	TRP	3.5
1	B	289	ASN	3.5
1	B	332	ASP	3.5
1	B	295	SER	3.4
1	B	319	GLU	3.4
1	A	564	PRO	3.4
1	B	453	ALA	3.4
1	B	487	GLY	3.4
1	A	563	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	494	ILE	3.4
1	A	267	VAL	3.4
1	B	76	VAL	3.4
1	B	49	GLY	3.4
1	B	77	VAL	3.3
1	B	326	PRO	3.3
1	B	308	THR	3.3
1	B	58	LEU	3.3
1	B	281	HIS	3.3
1	B	335	CYS	3.3
1	B	430	THR	3.3
1	A	263	ALA	3.2
1	B	287	VAL	3.3
1	B	246	ASP	3.2
1	B	28	ILE	3.2
1	B	450	SER	3.2
1	A	309	ASN	3.2
1	B	285	VAL	3.2
1	B	481	PRO	3.2
1	A	241	VAL	3.1
1	A	266	ILE	3.1
1	B	334	PHE	3.1
1	A	557	ALA	3.1
1	B	302	ASN	3.1
1	B	313	ASP	3.1
1	A	569	VAL	3.1
1	A	279	LEU	3.1
1	B	563	LEU	3.1
1	B	269	THR	3.1
1	B	297	THR	3.1
1	B	331	PRO	3.1
1	A	77	VAL	3.1
1	B	291	SER	3.0
1	A	256	HIS	3.0
1	B	270	GLN	3.0
1	B	57	GLU	3.0
1	B	52	ILE	2.9
1	B	359	TYR	2.9
1	B	258	GLY	2.9
1	B	278	ASP	2.9
1	B	484	VAL	2.9
1	A	271	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	461	VAL	2.9
1	B	490	ASN	2.9
1	B	55	LEU	2.9
1	B	328	PHE	2.9
1	A	254	VAL	2.9
1	B	426	ILE	2.9
1	A	568	ALA	2.8
1	B	492	ALA	2.8
1	B	243	PHE	2.8
1	B	263	ALA	2.8
1	B	316	LYS	2.8
1	A	571	THR	2.8
1	B	330	HIS	2.8
1	B	41	LEU	2.8
1	A	59	PRO	2.7
1	B	323	LYS	2.7
1	B	274	ASP	2.7
1	B	299	GLY	2.7
1	B	24	SER	2.7
1	A	409	TRP	2.7
1	B	353	ASN	2.7
1	B	240	GLY	2.6
1	B	32	LEU	2.6
1	B	233	LEU	2.6
1	B	565	GLU	2.6
1	B	250	LYS	2.6
1	B	25	ALA	2.6
1	B	236	LEU	2.6
1	A	558	ASP	2.6
1	B	26	TYR	2.6
1	B	234	ALA	2.5
1	B	312	GLY	2.5
1	B	264	LYS	2.5
1	B	311	LYS	2.5
1	B	59	PRO	2.5
1	B	248	HIS	2.5
1	B	293	THR	2.5
1	A	50	GLU	2.5
1	B	322	ALA	2.5
1	A	19	GLY	2.5
1	B	462	PRO	2.5
1	A	249	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	483	VAL	2.4
1	A	272	GLY	2.4
1	B	336	GLU	2.4
1	B	280	THR	2.4
1	B	350	THR	2.4
1	B	468	ILE	2.4
1	A	555	LYS	2.4
1	B	35	LEU	2.4
1	A	298	TYR	2.4
1	B	337	ASN	2.4
1	B	346	SER	2.4
1	A	581	VAL	2.4
1	B	434	ILE	2.4
1	B	464	TYR	2.4
1	B	290	GLY	2.4
1	A	565	GLU	2.4
1	B	38	ALA	2.3
1	A	591	VAL	2.3
1	A	587	GLU	2.3
1	B	404	ASP	2.3
1	B	498	ALA	2.3
1	A	280	THR	2.3
1	B	344	PHE	2.3
1	A	567	LEU	2.3
1	B	429	CYS	2.3
1	B	231	PRO	2.3
1	B	470	SER	2.3
1	B	294	GLU	2.3
1	B	34	GLY	2.3
1	B	225	TYR	2.3
1	B	409	TRP	2.3
1	B	283	ASP	2.2
1	A	26	TYR	2.2
1	A	53	HIS	2.2
1	A	269	THR	2.2
1	B	460	THR	2.2
1	B	486	GLU	2.2
1	B	454	SER	2.2
1	A	577	LYS	2.2
1	B	273	LYS	2.2
1	B	20	VAL	2.2
1	B	100	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	251	ASN	2.2
1	B	80	GLY	2.2
1	B	275	LYS	2.2
1	A	78	ARG	2.2
1	A	321	LEU	2.2
1	A	325	ASP	2.2
1	B	463	VAL	2.2
1	B	476	ARG	2.1
1	B	523	LEU	2.1
1	B	56	GLU	2.1
1	B	103	PRO	2.1
1	A	322	ALA	2.1
1	B	50	GLU	2.1
1	B	45	ALA	2.1
1	B	351	LEU	2.1
1	B	482	ASP	2.1
1	B	571	THR	2.1
1	A	41	LEU	2.1
1	B	256	HIS	2.1
1	B	452	LEU	2.1
1	B	524	LEU	2.1
1	B	284	ILE	2.1
1	B	428	ASP	2.1
1	B	431	GLY	2.1
1	A	274	ASP	2.0
1	A	278	ASP	2.0
1	B	402	ASN	2.0
1	B	466	PRO	2.0
1	B	40	PHE	2.0
1	B	48	LYS	2.0
1	B	265	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 oligosaccharides in this entry.

6.4 Ligands

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
7	GOL	B	608	6/6	0.66	0.22	73,84,90,97	0
4	MPD	B	604	8/8	0.79	0.21	39,45,54,60	0
7	GOL	B	606	6/6	0.79	0.18	48,59,61,65	0
4	MPD	A	606	8/8	0.79	0.27	75,84,87,87	0
5	NA	B	601	1/1	0.81	0.17	58,58,58,58	0
7	GOL	B	607	6/6	0.85	0.26	87,91,92,96	0
2	MES	B	602	12/12	0.86	0.15	49,67,81,84	0
6	PO4	B	603	5/5	0.88	0.13	50,67,70,83	0
2	MES	A	601	12/12	0.88	0.17	45,70,84,84	0
4	MPD	A	604	8/8	0.89	0.22	57,68,74,74	0
4	MPD	A	605	8/8	0.91	0.15	47,60,66,69	0
4	MPD	B	605	8/8	0.93	0.17	61,64,66,69	0
3	K	A	602	1/1	0.93	0.25	65,65,65,65	0
3	K	A	603	1/1	0.95	0.07	54,54,54,54	0

6.5 Other polymers

There are no such residues in this entry.