



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:19 PM GMT

PDB ID : 1T0S
Title : Structure of the Toluene/o-Xylene Monooxygenase Hydroxylase with 4-bromophenol bound
Authors : Sazinsky, M.H.; Bard, J.; Di Donato, A.; Lippard, S.J.
Deposited on : 2004-04-12
Resolution : 2.20 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

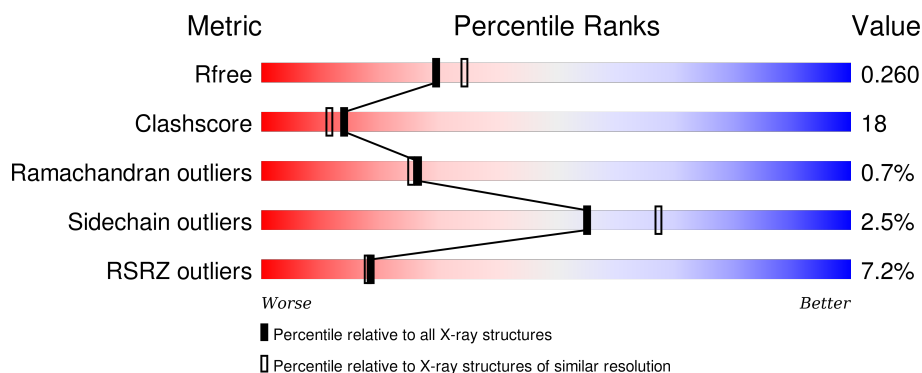
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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 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	498	<div> <div>6%</div> <div>63%</div> <div>34%</div> <div>..</div> </div>
2	B	330	<div> <div>4%</div> <div>76%</div> <div>19%</div> <div>..</div> </div>
3	C	86	<div> <div>23%</div> <div>42%</div> <div>52%</div> <div>..</div> </div>

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	BML	A	503	-	X	-	X
7	BML	A	504	-	X	-	-
7	BML	A	505	-	X	-	-
7	BML	A	506	-	X	-	X
7	BML	B	331	-	X	-	-
7	BML	B	332	-	X	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7667 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 toluene, o-xylene monooxygenase oxygenase subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	491	Total	C	N	O	S	0	0	0
			4018	2565	674	753	26			

- Molecule 2 is a protein called toluene, o-xylene monooxygenase oxygenase subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	323	Total	C	N	O	S	0	0	0
			2650	1680	468	492	10			

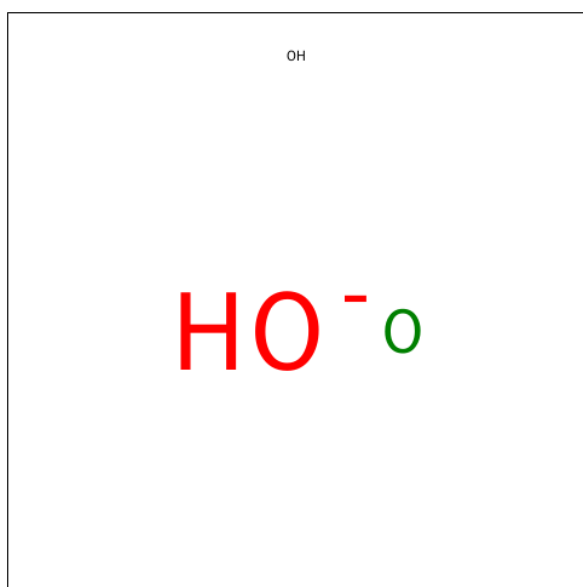
- Molecule 3 is a protein called touB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	83	Total	C	N	O	S	0	0	0
			676	425	120	126	5			

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

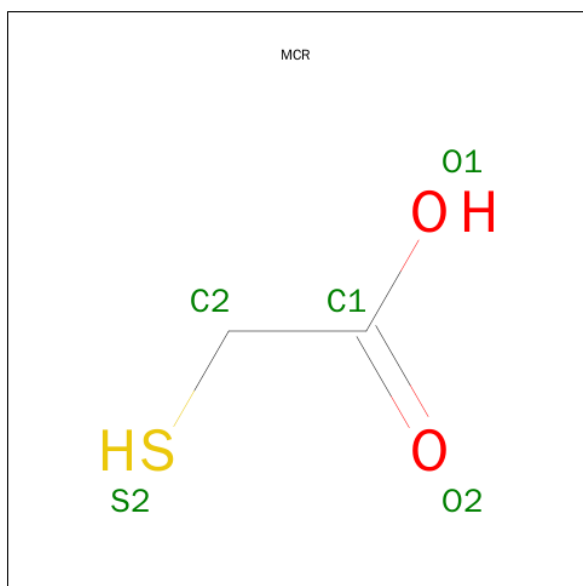
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Fe	0	0
			2	2		

- Molecule 5 is HYDROXIDE ION (three-letter code: OH) (formula: HO).



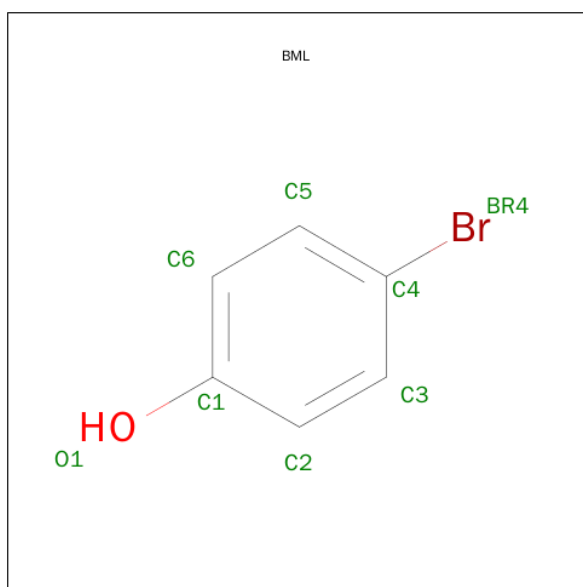
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O		0	0
			1	1			

- Molecule 6 is SULFANYLACETIC ACID (three-letter code: MCR) (formula: C₂H₄O₂S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	O	S	0	0
			5	2	2	1		

- Molecule 7 is 4-BROMOPHENOL (three-letter code: BML) (formula: C₆H₅BrO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	Br	C	O	0	0
			8	1	6	1		
7	A	1	Total	Br	C	O	0	0
			8	1	6	1		
7	A	1	Total	Br	C	O	0	0
			8	1	6	1		
7	B	1	Total	Br	C	O	0	0
			8	1	6	1		
7	A	1	Total	Br	C	O	0	0
			8	1	6	1		
7	B	1	Total	Br	C	O	0	0
			8	1	6	1		

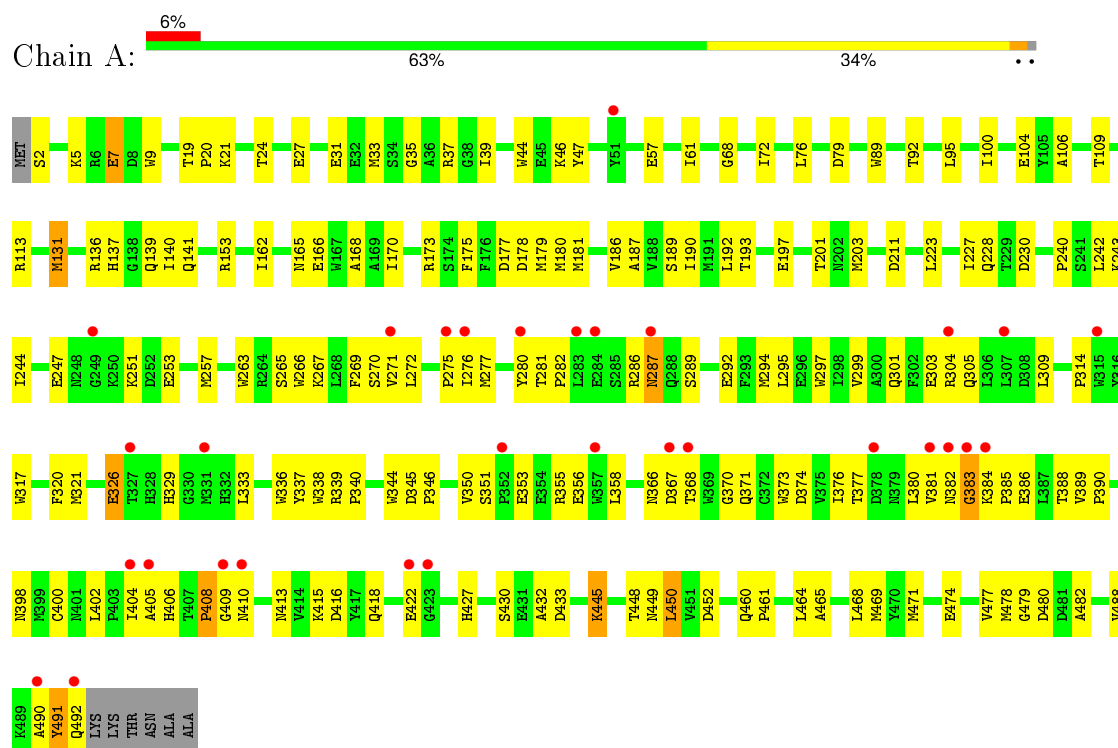
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	123	Total	O	0	0
			123	123		
8	B	141	Total	O	0	0
			141	141		
8	C	3	Total	O	0	0
			3	3		

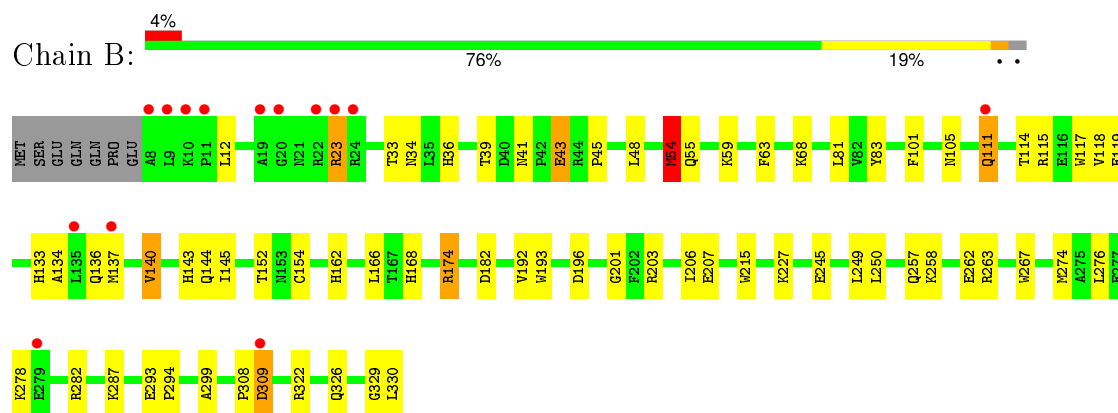
3 Residue-property plots

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.

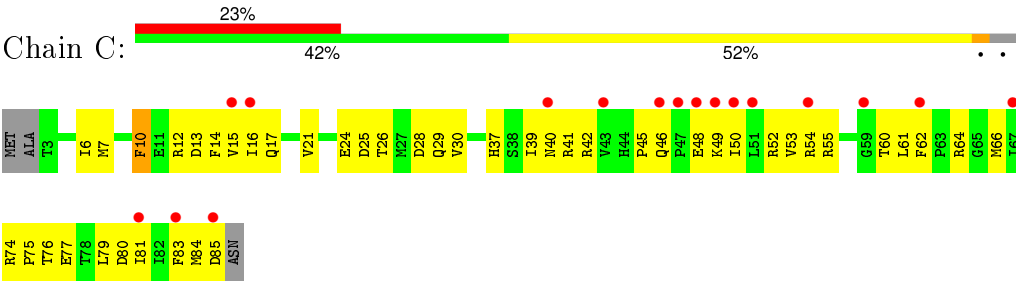
- Molecule 1: toluene, o-xylene monooxygenase oxygenase subunit



- Molecule 2: toluene, o-xylene monooxygenase oxygenase subunit



- Molecule 3: touB



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	183.24Å 183.24Å 67.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.47 – 2.20 29.47 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.5 (29.47-2.20) 91.3 (29.47-2.10)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.96 (at 2.10Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.236 , 0.273 0.230 , 0.260	Depositor DCC
R_{free} test set	2301 reflections (4.48%)	DCC
Wilson B-factor (Å ²)	37.1	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 67.0	EDS
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 72497 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7667	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.85% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FE, MCR, BML, OH

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	0.39	1/4142 (0.0%)	0.57	0/5629
2	B	0.45	2/2722 (0.1%)	0.60	0/3699
3	C	0.30	0/690	0.56	0/934
All	All	0.41	3/7554 (0.0%)	0.58	0/10262

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	119	GLU	CD-OE2	7.24	1.33	1.25
1	A	131	MET	SD-CE	-5.30	1.48	1.77
2	B	54	MET	SD-CE	-5.28	1.48	1.77

There are no bond angle outliers.

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	4018	0	3773	162	0
2	B	2650	0	2548	76	0
3	C	676	0	667	55	0
4	A	2	0	0	0	0
5	A	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	5	0	3	0	0
7	A	32	0	16	3	0
7	B	16	0	8	1	0
8	A	123	0	0	3	0
8	B	141	0	0	9	0
8	C	3	0	0	0	0
All	All	7667	0	7015	264	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 18.

All (264) 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:2:SER:N	2:B:105:ASN:HD22	1.76	0.84
1:A:139:GLN:HE22	2:B:83:TYR:H	1.25	0.82
1:A:131:MET:HE1	2:B:140:VAL:HG23	1.60	0.81
3:C:13:ASP:OD1	3:C:42:ARG:HD2	1.84	0.77
2:B:54:MET:HA	2:B:54:MET:HE2	1.66	0.77
1:A:106:ALA:HB1	1:A:181:MET:HE3	1.69	0.74
2:B:168:HIS:HD2	2:B:257:GLN:HE21	1.35	0.74
2:B:111:GLN:H	2:B:111:GLN:HE21	1.37	0.71
1:A:366:ASN:HA	1:A:370:GLY:HA3	1.71	0.71
1:A:140:ILE:HG21	1:A:227:ILE:HD11	1.72	0.71
1:A:7:GLU:H	1:A:7:GLU:CD	1.93	0.71
1:A:170:ILE:HD13	1:A:173:ARG:NH2	2.06	0.71
2:B:309:ASP:HB3	8:B:426:HOH:O	1.91	0.70
1:A:9:TRP:HB3	2:B:174:ARG:HG3	1.73	0.70
1:A:68:GLY:O	1:A:72:ILE:HG12	1.90	0.70
3:C:66:MET:CE	3:C:71:ALA:HB2	2.22	0.70
1:A:131:MET:CE	2:B:140:VAL:HG23	2.23	0.69
1:A:377:THR:O	1:A:381:VAL:HG23	1.92	0.69
1:A:276:ILE:O	1:A:281:THR:HG23	1.94	0.67
2:B:329:GLY:O	2:B:330:LEU:HB2	1.95	0.66
3:C:10:PHE:CE1	3:C:81:ILE:HG21	2.31	0.66
1:A:353:GLU:O	1:A:356:GLU:HG2	1.95	0.66
1:A:377:THR:HG21	1:A:491:TYR:CD1	2.31	0.66
1:A:367:ASP:HB3	1:A:410:ASN:HD22	1.62	0.65
1:A:405:ALA:HB1	3:C:14:PHE:O	1.97	0.64
1:A:405:ALA:HB2	3:C:15:VAL:HB	1.78	0.64
2:B:162:HIS:HE1	2:B:227:LYS:NZ	1.95	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:ASN:ND2	1:A:449:ASN:ND2	2.45	0.64
2:B:168:HIS:CD2	2:B:257:GLN:HE21	2.16	0.64
1:A:240:PRO:O	1:A:244:ILE:HG12	1.98	0.64
1:A:314:PRO:HD2	1:A:317:TRP:CE3	2.33	0.64
1:A:39:ILE:HB	1:A:44:TRP:NE1	2.13	0.64
2:B:162:HIS:HE1	2:B:227:LYS:HZ2	1.45	0.63
1:A:464:LEU:HD23	1:A:465:ALA:N	2.13	0.63
1:A:491:TYR:O	1:A:492:GLN:HB2	1.99	0.63
2:B:192:VAL:HG13	8:B:432:HOH:O	1.98	0.63
2:B:182:ASP:HB3	8:B:462:HOH:O	1.97	0.62
1:A:113:ARG:HH11	2:B:144:GLN:HE21	1.48	0.62
1:A:295:LEU:HD23	1:A:299:VAL:HG21	1.82	0.61
3:C:66:MET:HE3	3:C:71:ALA:HB2	1.83	0.61
2:B:258:LYS:O	2:B:262:GLU:HG3	2.00	0.61
1:A:382:ASN:O	1:A:384:LYS:HG3	2.00	0.61
2:B:33:THR:HB	2:B:34:ASN:HD22	1.65	0.61
2:B:41:ASN:HB3	2:B:43:GLU:OE2	2.00	0.60
2:B:111:GLN:H	2:B:111:GLN:NE2	1.99	0.60
2:B:196:ASP:HB2	8:B:432:HOH:O	1.99	0.60
1:A:460:GLN:HA	1:A:461:PRO:C	2.21	0.60
2:B:322:ARG:O	2:B:326:GLN:HG3	2.02	0.60
2:B:134:ALA:CB	2:B:206:ILE:HD12	2.32	0.60
1:A:468:LEU:HD22	1:A:478:MET:HE3	1.83	0.60
1:A:289:SER:OG	1:A:292:GLU:HG3	2.02	0.59
1:A:189:SER:O	1:A:193:THR:OG1	2.19	0.59
3:C:28:ASP:OD2	3:C:64:ARG:HB3	2.02	0.59
1:A:57:GLU:O	1:A:61:ILE:HG12	2.02	0.59
1:A:398:ASN:ND2	1:A:449:ASN:HD22	2.00	0.59
1:A:344:TRP:O	1:A:346:PRO:HD3	2.02	0.59
1:A:384:LYS:N	1:A:385:PRO:HD3	2.17	0.58
1:A:95:LEU:HD23	1:A:276:ILE:HD11	1.84	0.58
1:A:422:GLU:OE1	1:A:445:LYS:HE3	2.04	0.57
1:A:137:HIS:CD2	1:A:227:ILE:HG23	2.39	0.57
1:A:336:TRP:O	1:A:339:ARG:HB3	2.04	0.57
1:A:271:VAL:HG13	1:A:272:LEU:HG	1.86	0.57
1:A:253:GLU:CD	1:A:253:GLU:H	2.09	0.56
1:A:46:LYS:HD2	8:A:617:HOH:O	2.03	0.56
1:A:416:ASP:HB2	3:C:16:ILE:CD1	2.36	0.56
1:A:265:SER:O	1:A:269:PHE:HB2	2.04	0.56
1:A:37:ARG:NH2	1:A:257:MET:HE1	2.21	0.56
2:B:287:LYS:HB2	2:B:287:LYS:NZ	2.21	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:GLY:C	1:A:385:PRO:HD3	2.26	0.56
3:C:74:ARG:O	3:C:77:GLU:HB2	2.06	0.56
1:A:197:GLU:HA	1:A:201:THR:OG1	2.06	0.56
1:A:339:ARG:HH12	1:A:389:VAL:CG1	2.19	0.55
3:C:62:PHE:HB3	3:C:66:MET:HE2	1.88	0.55
1:A:39:ILE:HB	1:A:44:TRP:HE1	1.72	0.55
1:A:345:ASP:HB3	1:A:482:ALA:HA	1.89	0.55
1:A:47:TYR:CE1	1:A:240:PRO:HB2	2.42	0.55
3:C:54:ARG:HB3	3:C:61:LEU:HA	1.89	0.55
2:B:274:MET:HE2	8:B:420:HOH:O	2.07	0.54
2:B:133:HIS:O	2:B:136:GLN:HB3	2.07	0.54
2:B:43:GLU:CD	2:B:43:GLU:H	2.09	0.54
2:B:114:THR:O	2:B:118:VAL:HG23	2.07	0.54
1:A:413:ASN:O	1:A:415:LYS:HG3	2.06	0.54
1:A:136:ARG:NH1	1:A:230:ASP:OD2	2.40	0.54
1:A:281:THR:HG22	7:A:505:BML:C4	2.38	0.54
1:A:251:LYS:NZ	1:A:251:LYS:HB3	2.23	0.54
2:B:276:LEU:HD22	2:B:282:ARG:HB2	1.89	0.54
1:A:33:MET:SD	2:B:54:MET:HE1	2.48	0.54
1:A:427:HIS:CE1	3:C:76:THR:HG22	2.43	0.53
1:A:371:GLN:O	1:A:374:ASP:HB2	2.08	0.53
1:A:277:MET:HA	1:A:281:THR:CG2	2.39	0.53
3:C:6:ILE:HD12	3:C:6:ILE:C	2.28	0.53
3:C:53:VAL:HG22	3:C:81:ILE:HG13	1.90	0.53
1:A:180:MET:HB3	1:A:192:LEU:HD22	1.89	0.53
1:A:131:MET:CE	2:B:140:VAL:CG2	2.87	0.53
3:C:73:LEU:HD22	3:C:77:GLU:HG2	1.91	0.53
3:C:46:GLN:HB2	3:C:83:PHE:CE1	2.43	0.53
2:B:250:LEU:HG	8:B:440:HOH:O	2.08	0.53
1:A:474:GLU:O	1:A:478:MET:HG3	2.09	0.53
1:A:338:TRP:CD1	1:A:390:PRO:HG3	2.43	0.53
1:A:37:ARG:HH21	1:A:257:MET:CE	2.22	0.53
1:A:416:ASP:OD2	1:A:418:GLN:HG3	2.09	0.52
1:A:339:ARG:HG2	1:A:340:PRO:HD3	1.90	0.52
1:A:409:GLY:O	3:C:12:ARG:HG2	2.09	0.52
3:C:66:MET:HE1	3:C:71:ALA:HB2	1.91	0.52
1:A:211:ASP:OD2	7:A:504:BML:H6	2.09	0.52
2:B:193:TRP:CZ2	2:B:206:ILE:HD11	2.46	0.51
3:C:39:ILE:CG2	3:C:45:PRO:HG3	2.40	0.51
3:C:75:PRO:O	3:C:76:THR:CB	2.57	0.51
1:A:282:PRO:O	1:A:286:ARG:HG3	2.11	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:SER:O	1:A:355:ARG:HG3	2.10	0.51
3:C:52:ARG:HG3	3:C:52:ARG:HH11	1.76	0.50
3:C:80:ASP:C	3:C:81:ILE:HD12	2.31	0.50
1:A:491:TYR:N	1:A:491:TYR:CD2	2.79	0.50
1:A:166:GLU:HA	1:A:471:MET:HB3	1.92	0.50
1:A:338:TRP:C	1:A:340:PRO:HD2	2.32	0.50
1:A:24:THR:OG1	1:A:27:GLU:HG3	2.12	0.49
1:A:270:SER:O	1:A:275:PRO:HD3	2.12	0.49
1:A:384:LYS:HA	1:A:386:GLU:OE2	2.11	0.49
2:B:134:ALA:HB1	2:B:206:ILE:HD12	1.93	0.49
2:B:39:THR:HG22	2:B:63:PHE:HE1	1.78	0.49
1:A:277:MET:HA	1:A:281:THR:HG21	1.94	0.49
1:A:190:ILE:HD12	1:A:242:LEU:CD2	2.43	0.49
2:B:101:PHE:HE2	2:B:249:LEU:HD21	1.78	0.49
8:A:623:HOH:O	3:C:13:ASP:HB3	2.11	0.49
3:C:62:PHE:CD1	3:C:66:MET:HE1	2.47	0.49
1:A:491:TYR:N	1:A:491:TYR:HD2	2.10	0.49
1:A:153:ARG:NH1	2:B:12:LEU:HD21	2.27	0.49
1:A:263:TRP:CE2	1:A:432:ALA:HB3	2.48	0.48
1:A:400:CYS:O	1:A:402:LEU:HG	2.12	0.48
1:A:2:SER:N	2:B:105:ASN:ND2	2.54	0.48
2:B:111:GLN:N	2:B:111:GLN:NE2	2.62	0.48
1:A:295:LEU:HA	1:A:299:VAL:CG2	2.43	0.48
3:C:7:MET:HE2	3:C:76:THR:HG23	1.95	0.48
1:A:76:LEU:O	1:A:79:ASP:HB3	2.14	0.48
3:C:7:MET:CE	3:C:76:THR:HG23	2.44	0.48
2:B:117:TRP:CE3	2:B:245:GLU:HG3	2.49	0.48
1:A:267:LYS:NZ	1:A:433:ASP:OD2	2.44	0.48
1:A:243:LYS:O	1:A:247:GLU:HB2	2.14	0.48
3:C:55:ARG:HA	3:C:79:LEU:HD23	1.95	0.48
1:A:175:PHE:CE1	1:A:179:MET:HG3	2.48	0.48
3:C:10:PHE:HE2	3:C:39:ILE:HD11	1.78	0.48
1:A:339:ARG:NH1	1:A:389:VAL:HG12	2.28	0.48
1:A:153:ARG:CZ	2:B:12:LEU:HD21	2.43	0.48
3:C:81:ILE:N	3:C:81:ILE:HD12	2.29	0.47
1:A:377:THR:HG21	1:A:491:TYR:CE1	2.49	0.47
3:C:39:ILE:HG22	3:C:40:ASN:ND2	2.28	0.47
1:A:353:GLU:CD	1:A:353:GLU:H	2.18	0.47
2:B:192:VAL:HA	8:B:432:HOH:O	2.14	0.47
1:A:339:ARG:O	1:A:479:GLY:HA3	2.14	0.47
1:A:37:ARG:HH21	1:A:257:MET:HE3	1.79	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:355:ARG:HA	1:A:358:LEU:HD12	1.96	0.47
1:A:404:ILE:O	3:C:15:VAL:HG23	2.15	0.47
1:A:337:TYR:CE2	3:C:42:ARG:NH2	2.82	0.47
1:A:106:ALA:CB	1:A:181:MET:HE3	2.41	0.47
3:C:39:ILE:HG23	3:C:45:PRO:HG3	1.96	0.47
1:A:367:ASP:HB3	1:A:410:ASN:ND2	2.28	0.47
1:A:317:TRP:O	1:A:321:MET:HG2	2.15	0.47
1:A:37:ARG:HH11	1:A:37:ARG:HG3	1.79	0.47
1:A:294:MET:O	1:A:299:VAL:HG23	2.15	0.47
1:A:37:ARG:NH2	1:A:257:MET:CE	2.78	0.47
3:C:25:ASP:HB3	3:C:29:GLN:HB2	1.97	0.47
1:A:162:ILE:HG22	1:A:162:ILE:O	2.15	0.46
2:B:201:GLY:HA3	2:B:299:ALA:HA	1.97	0.46
2:B:154:CYS:HB3	2:B:267:TRP:CE2	2.50	0.46
1:A:400:CYS:SG	1:A:402:LEU:HB2	2.56	0.46
1:A:339:ARG:CD	1:A:480:ASP:HA	2.45	0.46
2:B:206:ILE:HG13	2:B:207:GLU:N	2.31	0.45
1:A:186:VAL:O	1:A:190:ILE:HG12	2.16	0.45
2:B:54:MET:HA	2:B:54:MET:CE	2.41	0.45
1:A:450:LEU:HA	1:A:450:LEU:HD12	1.74	0.45
1:A:165:ASN:OD1	1:A:170:ILE:HD11	2.17	0.45
2:B:23:ARG:O	2:B:23:ARG:HD2	2.17	0.45
1:A:287:ASN:N	1:A:287:ASN:HD22	2.14	0.45
1:A:104:GLU:HG3	1:A:141:GLN:NE2	2.32	0.45
1:A:309:LEU:N	1:A:309:LEU:HD12	2.31	0.45
2:B:81:LEU:HD11	2:B:263:ARG:HD3	1.98	0.45
1:A:177:ASP:HA	1:A:181:MET:HB2	1.98	0.45
1:A:281:THR:HG22	7:A:505:BML:BR4	2.72	0.45
1:A:405:ALA:CB	3:C:15:VAL:HB	2.46	0.45
1:A:474:GLU:O	1:A:477:VAL:HG22	2.16	0.45
1:A:350:VAL:HG11	1:A:373:TRP:CD2	2.51	0.45
3:C:75:PRO:O	3:C:76:THR:HB	2.17	0.45
2:B:45:PRO:HG2	2:B:55:GLN:OE1	2.17	0.45
3:C:12:ARG:HH11	3:C:12:ARG:HG3	1.81	0.44
2:B:12:LEU:N	2:B:12:LEU:HD22	2.33	0.44
1:A:301:GLN:H	1:A:301:GLN:CD	2.20	0.44
1:A:2:SER:HB2	2:B:105:ASN:HD21	1.83	0.44
3:C:41:ARG:HH11	3:C:41:ARG:HG2	1.82	0.44
1:A:137:HIS:CD2	1:A:227:ILE:HD12	2.52	0.44
3:C:17:GLN:NE2	3:C:37:HIS:HB3	2.33	0.44
1:A:21:LYS:HB3	2:B:203:ARG:NH2	2.32	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:VAL:O	1:A:303:GLU:HG3	2.18	0.44
2:B:143:HIS:CD2	2:B:143:HIS:C	2.91	0.44
1:A:136:ARG:HB2	2:B:83:TYR:CE2	2.51	0.44
1:A:113:ARG:HH11	2:B:144:GLN:NE2	2.14	0.44
1:A:326:GLU:O	1:A:329:HIS:HB2	2.17	0.44
2:B:68:LYS:HB2	2:B:278:LYS:HE2	2.00	0.44
3:C:53:VAL:O	3:C:62:PHE:HB2	2.17	0.44
1:A:287:ASN:ND2	1:A:287:ASN:N	2.66	0.44
2:B:308:PRO:O	2:B:309:ASP:HB2	2.17	0.43
3:C:25:ASP:HA	3:C:29:GLN:OE1	2.17	0.43
1:A:203:MET:HG2	1:A:297:TRP:HB3	1.99	0.43
1:A:95:LEU:HD13	1:A:344:TRP:CZ3	2.53	0.43
3:C:54:ARG:HB2	3:C:60:THR:O	2.18	0.43
1:A:131:MET:HE1	2:B:140:VAL:CG2	2.39	0.43
3:C:7:MET:HB3	3:C:16:ILE:CG2	2.49	0.43
1:A:263:TRP:O	1:A:266:TRP:HB3	2.19	0.43
3:C:49:LYS:C	3:C:50:ILE:HD12	2.38	0.43
1:A:340:PRO:HA	1:A:479:GLY:O	2.19	0.43
2:B:162:HIS:CE1	2:B:227:LYS:NZ	2.82	0.43
2:B:174:ARG:HG2	2:B:174:ARG:NH2	2.34	0.43
1:A:178:ASP:HA	2:B:48:LEU:HD11	2.00	0.42
1:A:314:PRO:O	1:A:317:TRP:HB2	2.19	0.42
1:A:468:LEU:HD13	1:A:478:MET:HE3	2.01	0.42
3:C:6:ILE:HD12	3:C:6:ILE:O	2.19	0.42
1:A:190:ILE:HD12	1:A:242:LEU:HG	2.00	0.42
1:A:166:GLU:OE1	1:A:168:ALA:HB3	2.20	0.42
1:A:31:GLU:OE2	1:A:35:GLY:HA2	2.20	0.42
1:A:380:LEU:HD21	1:A:388:THR:HG21	2.01	0.42
2:B:36:HIS:HE1	2:B:152:THR:OG1	2.01	0.42
1:A:106:ALA:O	1:A:181:MET:HE3	2.19	0.42
2:B:134:ALA:HA	2:B:137:MET:HE3	2.02	0.42
3:C:24:GLU:OE2	3:C:24:GLU:HA	2.20	0.42
1:A:368:THR:HA	1:A:409:GLY:HA3	2.02	0.42
2:B:59:LYS:HA	2:B:63:PHE:HD2	1.85	0.42
1:A:79:ASP:HB3	8:A:553:HOH:O	2.18	0.42
1:A:223:LEU:O	1:A:227:ILE:HG12	2.20	0.42
2:B:293:GLU:N	2:B:294:PRO:CD	2.83	0.42
1:A:448:THR:HB	1:A:452:ASP:HB2	2.02	0.42
3:C:21:VAL:HG11	3:C:30:VAL:HG13	2.01	0.42
1:A:333:LEU:HD23	3:C:14:PHE:CG	2.54	0.41
3:C:54:ARG:CB	3:C:61:LEU:HA	2.50	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:MET:CB	1:A:192:LEU:HD22	2.50	0.41
3:C:39:ILE:O	3:C:40:ASN:HB2	2.20	0.41
1:A:449:ASN:ND2	1:A:452:ASP:OD2	2.53	0.41
2:B:258:LYS:HE2	7:B:332:BML:BR4	2.76	0.41
2:B:145:ILE:HG23	2:B:215:TRP:HB2	2.02	0.41
3:C:17:GLN:HE22	3:C:37:HIS:HB3	1.85	0.41
3:C:52:ARG:NH1	3:C:52:ARG:HG3	2.35	0.41
3:C:84:MET:HB3	3:C:85:ASP:H	1.67	0.41
1:A:301:GLN:O	1:A:305:GLN:HG2	2.20	0.41
1:A:19:THR:HA	1:A:20:PRO:HD2	1.89	0.41
1:A:106:ALA:C	1:A:181:MET:HE3	2.41	0.41
1:A:109:THR:HG23	2:B:144:GLN:HB2	2.03	0.41
2:B:193:TRP:CZ2	2:B:203:ARG:HG3	2.56	0.41
3:C:7:MET:HE2	3:C:76:THR:CG2	2.51	0.41
1:A:2:SER:HB2	2:B:105:ASN:ND2	2.35	0.41
1:A:131:MET:HE3	2:B:140:VAL:CG2	2.51	0.41
2:B:154:CYS:HB3	2:B:267:TRP:CD2	2.56	0.41
1:A:5:LYS:HB3	1:A:7:GLU:OE1	2.20	0.41
2:B:192:VAL:HG22	8:B:432:HOH:O	2.21	0.41
1:A:468:LEU:HD13	1:A:478:MET:CE	2.51	0.41
2:B:287:LYS:HZ2	2:B:287:LYS:HB2	1.86	0.41
1:A:305:GLN:O	1:A:309:LEU:HD13	2.20	0.41
1:A:100:ILE:CG2	1:A:104:GLU:HG2	2.51	0.41
1:A:187:ALA:HB2	1:A:257:MET:HE1	2.02	0.40
2:B:145:ILE:O	2:B:145:ILE:HG13	2.21	0.40
1:A:280:TYR:CD2	1:A:280:TYR:N	2.89	0.40
1:A:266:TRP:CD2	1:A:320:PHE:HE1	2.39	0.40
3:C:26:THR:O	3:C:29:GLN:N	2.54	0.40
1:A:376:ILE:HG23	1:A:388:THR:HG22	2.02	0.40
1:A:488:VAL:C	1:A:490:ALA:N	2.74	0.40
1:A:89:TRP:O	1:A:92:THR:HB	2.20	0.40
2:B:115:ARG:HD3	8:B:472:HOH:O	2.21	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	489/498 (98%)	443 (91%)	41 (8%)	5 (1%)	19	16
2	B	321/330 (97%)	315 (98%)	6 (2%)	0	100	100
3	C	81/86 (94%)	70 (86%)	10 (12%)	1 (1%)	16	12
All	All	891/914 (98%)	828 (93%)	57 (6%)	6 (1%)	26	25

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	326	GLU
1	A	445	LYS
3	C	73	LEU
1	A	287	ASN
1	A	408	PRO
1	A	383	GLY

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	417/422 (99%)	408 (98%)	9 (2%)	60	72
2	B	275/282 (98%)	267 (97%)	8 (3%)	50	62
3	C	77/79 (98%)	75 (97%)	2 (3%)	54	66
All	All	769/783 (98%)	750 (98%)	19 (2%)	55	67

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLU
1	A	228	GLN
1	A	304	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	406	HIS
1	A	408	PRO
1	A	430	SER
1	A	450	LEU
1	A	469	MET
1	A	491	TYR
2	B	23	ARG
2	B	43	GLU
2	B	54	MET
2	B	111	GLN
2	B	140	VAL
2	B	166	LEU
2	B	174	ARG
2	B	309	ASP
3	C	10	PHE
3	C	48	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	GLN
1	A	204	GLN
1	A	228	GLN
1	A	237	GLN
1	A	248	ASN
1	A	287	ASN
1	A	322	GLN
1	A	379	ASN
1	A	398	ASN
1	A	410	ASN
1	A	418	GLN
1	A	427	HIS
2	B	34	ASN
2	B	36	HIS
2	B	87	ASN
2	B	98	GLN
2	B	111	GLN
2	B	144	GLN
2	B	153	ASN
2	B	162	HIS
2	B	168	HIS
2	B	286	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	326	GLN
3	C	40	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 1 is modelled with single atom and 2 are monoatomic - leaving 7 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
6	MCR	A	502	4	1,4,4	0.13	0	0,4,4	0.00	-
7	BML	A	503	-	8,8,8	3.41	5 (62%)	10,10,10	3.10	5 (50%)
7	BML	A	504	-	8,8,8	3.57	6 (75%)	10,10,10	2.95	5 (50%)
7	BML	A	505	-	8,8,8	3.50	6 (75%)	10,10,10	2.98	5 (50%)
7	BML	A	506	-	8,8,8	3.50	5 (62%)	10,10,10	3.08	5 (50%)
7	BML	B	331	-	8,8,8	3.53	5 (62%)	10,10,10	3.03	5 (50%)
7	BML	B	332	-	8,8,8	3.54	6 (75%)	10,10,10	2.99	5 (50%)

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
6	MCR	A	502	4	-	0/0/2/2	0/0/0/0
7	BML	A	503	-	-	0/0/0/0	0/1/1/1
7	BML	A	504	-	-	0/0/0/0	0/1/1/1
7	BML	A	505	-	-	0/0/0/0	0/1/1/1
7	BML	A	506	-	-	0/0/0/0	0/1/1/1
7	BML	B	331	-	-	0/0/0/0	0/1/1/1
7	BML	B	332	-	-	0/0/0/0	0/1/1/1

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	504	BML	C6-C5	-4.72	1.30	1.38
7	B	332	BML	C6-C5	-4.27	1.31	1.38
7	B	331	BML	C6-C5	-4.25	1.31	1.38
7	A	505	BML	C6-C5	-4.17	1.31	1.38
7	A	506	BML	C6-C5	-4.16	1.31	1.38
7	A	503	BML	C6-C5	-4.08	1.31	1.38
7	B	332	BML	BR4-C4	-2.30	1.85	1.90
7	A	504	BML	BR4-C4	-2.26	1.85	1.90
7	A	505	BML	BR4-C4	-2.01	1.85	1.90
7	A	503	BML	C3-C4	2.05	1.42	1.38
7	A	506	BML	C3-C4	2.17	1.43	1.38
7	A	505	BML	C3-C4	2.18	1.43	1.38
7	A	504	BML	C3-C4	2.21	1.43	1.38
7	B	332	BML	C3-C4	2.25	1.43	1.38
7	B	331	BML	C3-C4	2.38	1.43	1.38
7	A	504	BML	C6-C1	3.76	1.46	1.38
7	A	503	BML	C6-C1	3.89	1.46	1.38
7	A	505	BML	C6-C1	4.12	1.47	1.38
7	B	331	BML	C6-C1	4.15	1.47	1.38
7	B	332	BML	C6-C1	4.21	1.47	1.38
7	A	506	BML	C6-C1	4.23	1.47	1.38
7	B	332	BML	C3-C2	4.88	1.47	1.38
7	A	503	BML	C3-C2	4.88	1.47	1.38
7	A	505	BML	C3-C2	4.92	1.47	1.38
7	A	504	BML	C3-C2	4.96	1.47	1.38
7	A	506	BML	C3-C2	5.01	1.47	1.38
7	B	331	BML	C3-C2	5.05	1.47	1.38
7	B	332	BML	C2-C1	5.19	1.49	1.38
7	A	503	BML	C2-C1	5.25	1.49	1.38

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	505	BML	C2-C1	5.35	1.49	1.38
7	A	504	BML	C2-C1	5.36	1.49	1.38
7	A	506	BML	C2-C1	5.36	1.49	1.38
7	B	331	BML	C2-C1	5.40	1.49	1.38

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	506	BML	C3-C2-C1	-4.71	114.42	119.87
7	A	504	BML	C3-C2-C1	-4.70	114.43	119.87
7	B	332	BML	C3-C2-C1	-4.68	114.45	119.87
7	A	503	BML	C3-C2-C1	-4.67	114.47	119.87
7	B	331	BML	C3-C2-C1	-4.57	114.59	119.87
7	A	505	BML	C3-C2-C1	-4.55	114.61	119.87
7	A	506	BML	C5-C4-C3	-2.72	116.39	121.41
7	A	503	BML	C5-C4-C3	-2.71	116.42	121.41
7	B	331	BML	C5-C4-C3	-2.65	116.52	121.41
7	A	505	BML	C5-C4-C3	-2.51	116.78	121.41
7	A	504	BML	C5-C4-C3	-2.46	116.87	121.41
7	B	332	BML	C5-C4-C3	-2.44	116.91	121.41
7	A	504	BML	C2-C3-C4	2.19	122.34	119.17
7	A	505	BML	C2-C3-C4	2.38	122.62	119.17
7	B	331	BML	C2-C3-C4	2.40	122.65	119.17
7	B	332	BML	C2-C3-C4	2.40	122.65	119.17
7	A	503	BML	C2-C3-C4	2.63	122.99	119.17
7	A	506	BML	C2-C3-C4	2.69	123.07	119.17
7	A	504	BML	BR4-C4-C5	3.80	125.28	119.28
7	A	505	BML	BR4-C4-C5	4.03	125.65	119.28
7	B	332	BML	BR4-C4-C5	4.06	125.71	119.28
7	B	331	BML	BR4-C4-C5	4.09	125.75	119.28
7	A	506	BML	BR4-C4-C5	4.15	125.84	119.28
7	A	503	BML	BR4-C4-C5	4.32	126.11	119.28
7	B	332	BML	C6-C5-C4	5.73	127.49	119.17
7	A	505	BML	C6-C5-C4	5.83	127.63	119.17
7	A	504	BML	C6-C5-C4	5.88	127.70	119.17
7	A	503	BML	C6-C5-C4	5.90	127.73	119.17
7	A	506	BML	C6-C5-C4	5.91	127.75	119.17
7	B	331	BML	C6-C5-C4	6.00	127.88	119.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	504	BML	1	0
7	A	505	BML	2	0
7	B	332	BML	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	491/498 (98%)	0.26	31 (6%) 23 23	23, 49, 84, 103	0
2	B	323/330 (97%)	-0.11	14 (4%) 39 38	24, 35, 58, 90	0
3	C	83/86 (96%)	1.43	20 (24%) 1 1	44, 78, 94, 101	0
All	All	897/914 (98%)	0.24	65 (7%) 18 18	23, 43, 85, 103	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	381	VAL	5.5
3	C	59	GLY	5.2
3	C	72	GLY	4.7
2	B	22	ARG	4.5
3	C	47	PRO	4.4
3	C	50	ILE	4.1
1	A	492	GLN	4.1
3	C	48	GLU	4.0
1	A	423	GLY	3.9
3	C	83	PHE	3.9
2	B	8	ALA	3.9
1	A	51	TYR	3.9
3	C	15	VAL	3.7
1	A	383	GLY	3.5
1	A	410	ASN	3.5
3	C	16	ILE	3.3
1	A	490	ALA	3.2
1	A	352	PRO	3.1
3	C	85	ASP	3.1
1	A	304	ARG	3.1
2	B	20	GLY	3.0
1	A	382	ASN	3.0
2	B	10	LYS	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	23	ARG	3.0
2	B	9	LEU	2.9
3	C	68	VAL	2.8
1	A	368	THR	2.8
1	A	357	TRP	2.8
1	A	331	MET	2.8
1	A	280	TYR	2.8
1	A	287	ASN	2.7
3	C	46	GLN	2.7
3	C	40	ASN	2.7
2	B	111	GLN	2.7
1	A	271	VAL	2.7
3	C	62	PHE	2.6
2	B	135	LEU	2.6
3	C	43	VAL	2.5
1	A	284	GLU	2.5
1	A	276	ILE	2.5
3	C	51	LEU	2.4
1	A	249	GLY	2.4
3	C	73	LEU	2.4
1	A	367	ASP	2.3
1	A	378	ASP	2.3
1	A	404	ILE	2.3
1	A	275	PRO	2.3
2	B	24	ARG	2.3
1	A	307	LEU	2.3
1	A	315	TRP	2.3
2	B	11	PRO	2.2
1	A	409	GLY	2.2
2	B	309	ASP	2.2
1	A	384	LYS	2.2
3	C	81	ILE	2.2
3	C	54	ARG	2.1
1	A	327	THR	2.1
1	A	405	ALA	2.1
1	A	283	LEU	2.1
2	B	19	ALA	2.1
3	C	49	LYS	2.0
3	C	67	ILE	2.0
1	A	422	GLU	2.0
2	B	279	GLU	2.0
2	B	137	MET	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
7	BML	A	503	8/8	0.82	0.28	2.68	75,78,80,87	0
7	BML	A	506	8/8	0.92	0.23	2.27	71,73,75,79	0
7	BML	B	332	8/8	0.95	0.21	2.12	63,64,65,66	0
6	MCR	A	502	5/5	0.86	0.18	1.14	48,50,52,53	0
7	BML	A	504	8/8	0.88	0.20	0.33	72,73,75,81	0
7	BML	A	505	8/8	0.95	0.18	0.20	70,70,70,74	0
4	FE	A	500	1/1	0.98	0.09	-0.86	37,37,37,37	0
7	BML	B	331	8/8	0.97	0.10	-0.92	31,33,35,45	0
5	OH	A	501	1/1	1.00	0.06	-1.87	29,29,29,29	0
4	FE	A	499	1/1	0.99	0.05	-2.30	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.