



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:30 PM GMT

PDB ID : 1XL8  
Title : Crystal structure of mouse carnitine octanoyltransferase in complex with octanoylcarnitine  
Authors : Jogl, G.; Hsiao, Y.S.; Tong, L.  
Deposited on : 2004-09-30  
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](mailto: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.

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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

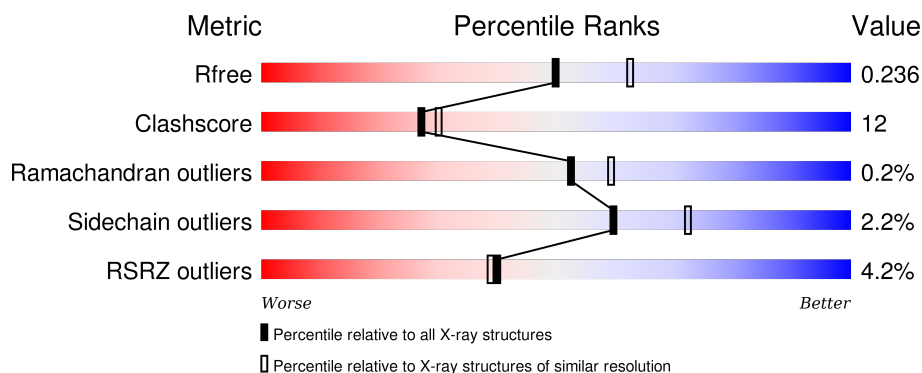
# 1 Overall quality at a glance

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  $\geq 3$ , 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	612	<div> <div>6%</div> <div> <div></div> <div>69%</div> <div>26%</div> <div>••</div> </div> </div>
1	B	612	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>20%</div> <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
2	152	A	613	-	-	-	X
4	MPD	B	614	-	-	-	X
4	MPD	B	615	-	-	-	X
4	MPD	B	616	-	-	-	X
4	MPD	B	618	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10118 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 Peroxisomal carnitine O-octanoyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	591	Total	C	N	O	S	Se	0	0	0
			4776	3042	829	872	16	17			
1	B	600	Total	C	N	O	S	Se	0	2	0
			4857	3093	841	888	17	18			

There are 34 discrepancies between the modelled and reference sequences:

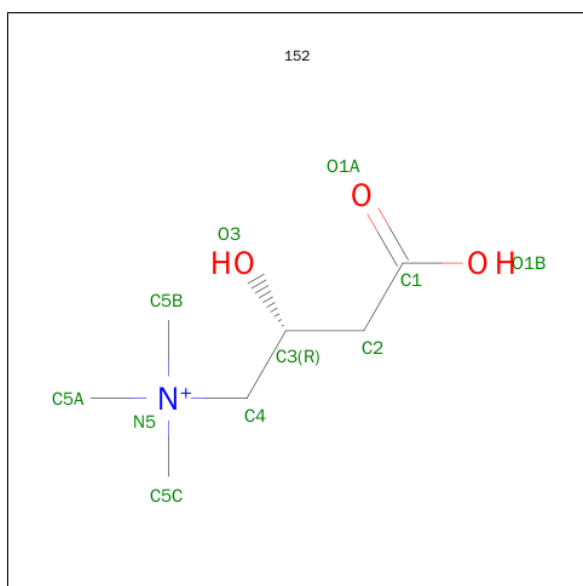
Chain	Residue	Modelled	Actual	Comment	Reference
A	127	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	128	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	156	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	161	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	177	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	296	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	333	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	335	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	443	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	469	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	483	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	493	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	494	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	558	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	595	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	603	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
A	607	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	127	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	128	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	156	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	161	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	177	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	296	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	333	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	335	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50

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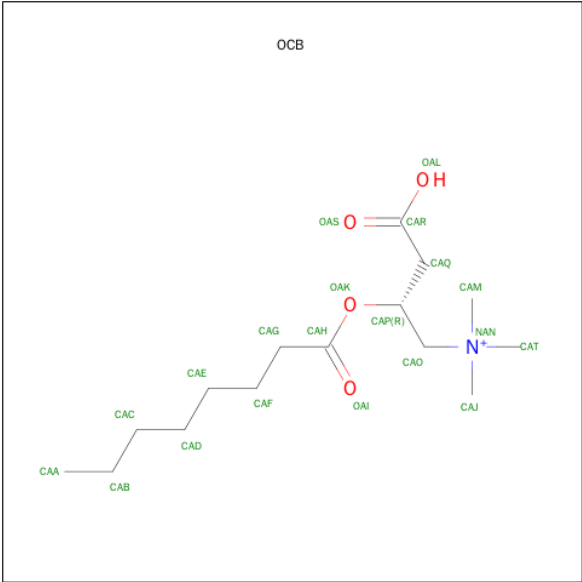
Chain	Residue	Modelled	Actual	Comment	Reference
B	443	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	469	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	483	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	493	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	494	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	558	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	595	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	603	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	607	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50

- Molecule 2 is CARNITINE (three-letter code: 152) (formula:  $C_7H_{16}NO_3$ ).



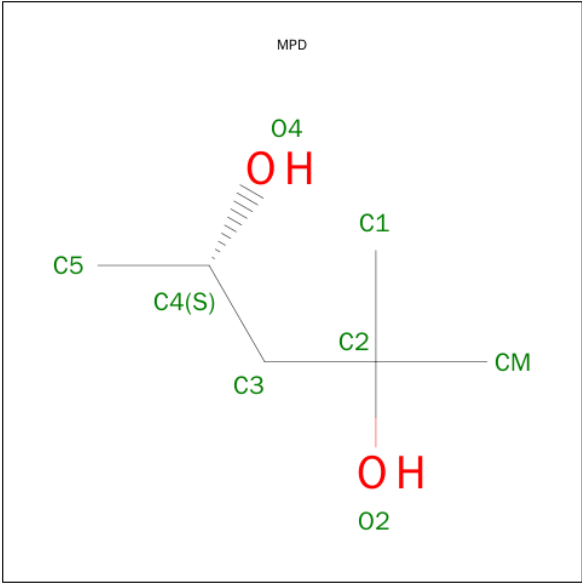
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	7	1	3		

- Molecule 3 is OCTANOYLCARNITINE (three-letter code: OCB) (formula:  $C_{15}H_{30}NO_4$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			20	15	1	4		

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>2</sub>).



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

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	138	Total	O	0	0
			138	138		
5	B	276	Total	O	0	0
			276	276		

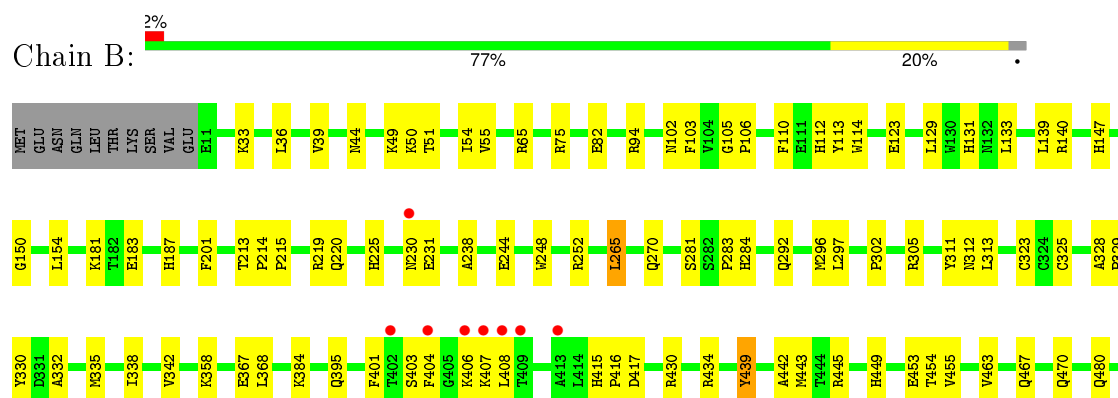
### 3 Residue-property plots

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.

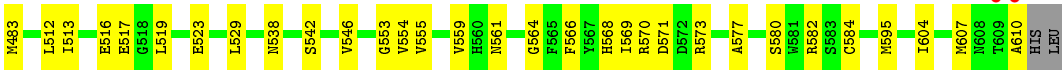
- Molecule 1: Peroxisomal carnitine O-octanoyltransferase



- Molecule 1: Peroxisomal carnitine O-octanoyltransferase







## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	163.61Å 163.61Å 159.48Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.09 – 2.20 29.08 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.2 (29.09-2.20) 95.4 (29.08-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.23 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.199 , 0.230 0.206 , 0.236	Depositor DCC
$R_{free}$ test set	5806 reflections (7.54%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.7	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 57.3	EDS
Estimated twinning fraction	0.019 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 80628 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10118	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MPD, 152, OCB

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	0.31	0/4884	0.54	0/6585
1	B	0.36	0/4966	0.60	2/6694 (0.0%)
All	All	0.34	0/9850	0.57	2/13279 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	564	GLY	N-CA-C	-6.41	97.06	113.10
1	B	150	GLY	N-CA-C	-5.29	99.86	113.10

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	4776	0	4676	123	0
1	B	4857	0	4762	105	0
2	A	11	0	15	3	0
3	B	20	0	29	8	0
4	B	40	0	70	2	0
5	A	138	0	0	4	0
5	B	276	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	10118	0	9552	231	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 12.

All (231) 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:568:HIS:HE1	1:A:570:ARG:HE	1.15	0.91
1:B:568:HIS:HE1	1:B:570:ARG:HE	1.17	0.89
1:A:219:ARG:HH21	1:A:220:GLN:HE22	1.20	0.89
1:B:252:ARG:HH21	1:B:270:GLN:HE22	1.23	0.86
1:B:480:GLN:HG3	1:B:607:MSE:HE1	1.58	0.84
1:B:430:ARG:HE	1:B:470:GLN:HE22	1.25	0.83
1:A:604:ILE:HA	1:A:607:MSE:HE2	1.60	0.81
1:A:483:MSE:HE3	1:A:606:LEU:HD23	1.64	0.80
1:B:284:HIS:H	1:B:292:GLN:NE2	1.79	0.80
1:A:430:ARG:HE	1:A:470:GLN:NE2	1.80	0.79
1:B:430:ARG:HE	1:B:470:GLN:NE2	1.83	0.77
1:A:430:ARG:HE	1:A:470:GLN:HE22	1.33	0.75
1:A:199:ARG:HH11	1:A:228:CYS:HB3	1.52	0.74
1:B:443:MSE:HE2	1:B:445:ARG:HD3	1.70	0.74
1:B:283:PRO:HA	1:B:292:GLN:HE21	1.53	0.73
1:A:345:ARG:HH21	1:A:348:GLU:CD	1.93	0.72
1:A:199:ARG:NH1	1:A:228:CYS:HB3	2.05	0.72
1:B:154:LEU:HD13	1:B:443:MSE:HE1	1.71	0.72
1:B:65:ARG:HH22	1:B:519:LEU:HD11	1.55	0.72
1:A:416:PRO:HB2	1:A:569:ILE:HD13	1.72	0.70
1:A:600:PHE:O	1:A:604:ILE:HG12	1.92	0.70
1:B:404:PHE:CZ	1:B:604:ILE:HG13	2.27	0.69
1:B:416:PRO:HB2	1:B:569:ILE:HD13	1.73	0.69
1:B:325:CYS:SG	1:B:335[B]:MSE:HE3	2.34	0.68
1:B:219:ARG:HH21	1:B:220:GLN:HE22	1.42	0.67
1:B:284:HIS:H	1:B:292:GLN:HE21	1.43	0.67
1:A:292:GLN:HE21	1:A:296:MSE:HE1	1.59	0.67
1:A:112:HIS:CD2	1:A:113:TYR:H	2.14	0.65
1:A:542:SER:OG	2:A:613:152:H5C1	1.97	0.64
1:B:570:ARG:HD2	1:B:573:ARG:HH11	1.62	0.64
1:B:358:LYS:NZ	1:B:358:LYS:HB3	2.13	0.63
1:B:415:HIS:HD2	1:B:417:ASP:H	1.45	0.63
1:B:335[A]:MSE:SE	3:B:613:OCB:HAA3	2.48	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:LEU:HD13	1:A:443:MSE:HE1	1.81	0.62
1:B:395:GLN:HG2	1:B:582:ARG:HG3	1.80	0.62
1:B:65:ARG:NH2	1:B:519:LEU:HD11	2.14	0.62
1:A:603:MSE:O	1:A:607:MSE:HG3	1.99	0.62
1:A:219:ARG:NH2	1:A:220:GLN:HE22	1.93	0.62
1:A:421:GLN:NE2	1:A:440:GLU:HG3	2.15	0.61
1:B:395:GLN:HG2	1:B:582:ARG:CG	2.31	0.61
1:B:147:HIS:CD2	1:B:445:ARG:HH22	2.19	0.61
1:A:263:LEU:O	1:A:267:GLU:HG3	2.02	0.60
1:B:51:THR:OG1	1:B:529:LEU:HD12	2.00	0.60
1:A:568:HIS:CE1	1:A:570:ARG:HE	2.08	0.59
1:B:54:ILE:HG23	1:B:523:GLU:HG3	1.84	0.59
1:A:121:GLN:OE1	1:A:278:ILE:HG23	2.03	0.59
1:B:302:PRO:HG2	1:B:384:LYS:HG3	1.83	0.59
1:A:292:GLN:HG2	1:A:296:MSE:HE3	1.86	0.58
1:B:415:HIS:CD2	1:B:417:ASP:H	2.21	0.58
1:A:483:MSE:CE	1:A:606:LEU:HD23	2.33	0.57
1:B:311:TYR:CE1	1:B:335[B]:MSE:SE	3.08	0.57
1:A:259:ASP:HB3	1:A:262:ASN:ND2	2.20	0.57
1:A:292:GLN:HE21	1:A:296:MSE:CE	2.18	0.56
1:A:69:LYS:HE3	1:A:517:GLU:OE1	2.05	0.56
1:A:465:TRP:CZ2	1:A:483:MSE:HE2	2.40	0.56
1:A:604:ILE:HA	1:A:607:MSE:CE	2.34	0.56
1:A:442:ALA:HB2	1:A:455:VAL:HG23	1.88	0.55
1:A:465:TRP:HZ2	1:A:483:MSE:HE2	1.71	0.55
1:A:608:ASN:O	1:A:611:HIS:HB2	2.07	0.55
1:B:33:LYS:NZ	5:B:775:HOH:O	2.40	0.55
1:A:109:HIS:HD2	1:A:343:ASP:OD1	1.90	0.55
1:B:407:LYS:CD	1:B:571:ASP:HB2	2.37	0.54
1:A:546:VAL:CG2	1:A:568:HIS:HB3	2.37	0.54
1:B:332:ALA:CB	3:B:613:OCB:HAD2	2.36	0.54
3:B:613:OCB:HAJ3	3:B:613:OCB:OAK	2.06	0.54
1:A:283:PRO:HA	1:A:292:GLN:HE21	1.72	0.54
2:A:613:152:H5B3	2:A:613:152:O3	2.08	0.54
1:A:36:LEU:HD21	1:A:52:GLU:HA	1.89	0.54
1:A:463:VAL:O	1:A:467:GLN:HG3	2.07	0.53
1:B:112:HIS:CD2	1:B:113:TYR:H	2.26	0.53
1:B:480:GLN:NE2	1:B:610:ALA:HB1	2.23	0.53
1:A:201:PHE:CE2	1:A:367:GLU:HB2	2.43	0.53
1:B:404:PHE:CE2	1:B:604:ILE:HG13	2.43	0.53
1:B:75:ARG:HG2	4:B:616:MPD:HM1	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:LEU:HD13	1:B:55:VAL:HG21	1.91	0.53
1:A:483:MSE:HE2	1:A:603:MSE:HG2	1.90	0.52
1:B:335[A]:MSE:CE	3:B:613:OCB:HAA3	2.39	0.52
1:B:554:VAL:CG1	1:B:577:ALA:HB1	2.40	0.52
1:A:546:VAL:HG22	1:A:567:TYR:O	2.09	0.52
1:A:52:GLU:O	1:A:56:GLN:HG3	2.10	0.52
1:B:512:LEU:O	1:B:516:GLU:HG3	2.10	0.52
1:A:591:LYS:O	1:A:595:MSE:HG2	2.10	0.52
1:A:242:SER:HB2	1:A:387:HIS:CD2	2.45	0.51
1:A:233:VAL:HG12	1:A:234:GLY:N	2.23	0.51
1:A:416:PRO:CB	1:A:569:ILE:HD13	2.41	0.51
1:A:68:GLN:O	1:A:72:GLU:HG3	2.10	0.51
1:A:51:THR:OG1	1:A:529:LEU:HD12	2.10	0.51
1:B:480:GLN:HE22	1:B:610:ALA:HB1	1.76	0.51
1:A:283:PRO:HA	1:A:292:GLN:NE2	2.25	0.51
1:B:404:PHE:CB	1:B:408:LEU:HD13	2.41	0.51
1:B:555:VAL:HG22	1:B:566:PHE:CZ	2.46	0.51
1:A:593:VAL:O	1:A:596:ILE:HG22	2.11	0.51
1:A:546:VAL:HG23	1:A:568:HIS:HB3	1.92	0.50
1:B:517:GLU:HB2	1:B:519:LEU:HG	1.93	0.50
1:B:36:LEU:O	1:B:39:VAL:HG22	2.10	0.50
1:A:43:ALA:HB2	1:A:529:LEU:HD21	1.94	0.50
1:A:560:HIS:HE1	1:A:583:SER:OG	1.93	0.50
1:A:252:ARG:O	1:A:256:ILE:HG13	2.12	0.49
1:B:442:ALA:HB2	1:B:455:VAL:HG23	1.94	0.49
1:B:546:VAL:HG12	1:B:546:VAL:O	2.13	0.49
1:A:12:ARG:O	1:A:15:GLN:HG2	2.12	0.49
1:A:219:ARG:NH2	1:A:362:ILE:O	2.38	0.49
1:B:554:VAL:CG1	1:B:577:ALA:CB	2.91	0.49
1:A:223:TYR:CD1	1:A:365:PRO:HD2	2.48	0.48
1:A:98:GLN:O	1:A:102:ASN:HB2	2.13	0.48
1:A:192:ILE:HD11	1:A:203:PHE:CZ	2.48	0.48
1:A:568:HIS:HE1	1:A:570:ARG:NE	1.97	0.47
1:B:463:VAL:O	1:B:467:GLN:HG3	2.13	0.47
1:B:607:MSE:HA	1:B:607:MSE:HE2	1.96	0.47
1:B:415:HIS:HD2	1:B:417:ASP:N	2.12	0.47
1:A:415:HIS:CD2	1:A:417:ASP:H	2.31	0.47
1:B:323[A]:CYS:HB2	1:B:335[A]:MSE:HE2	1.96	0.47
1:B:407:LYS:HD2	1:B:571:ASP:HB2	1.97	0.47
1:B:406:LYS:HG2	1:B:571:ASP:HA	1.95	0.47
1:B:439:TYR:CD1	1:B:439:TYR:C	2.88	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:ARG:HA	1:A:127:MSE:HE3	1.96	0.47
1:B:483:MSE:HE2	1:B:607:MSE:HE3	1.97	0.47
1:A:356:SER:OG	1:A:358:LYS:HG2	2.14	0.47
1:B:401:PHE:CE2	1:B:403:SER:HB2	2.49	0.47
1:A:249:ALA:O	1:A:253:GLU:HG3	2.13	0.47
1:B:110:PHE:HA	1:B:114:TRP:O	2.14	0.47
1:B:542:SER:OG	3:B:613:OCB:HAT1	2.14	0.46
1:A:252:ARG:HH21	1:A:270:GLN:HE22	1.63	0.46
1:A:293:VAL:O	1:A:297:LEU:HB2	2.16	0.46
1:A:36:LEU:O	1:A:39:VAL:HG22	2.15	0.46
1:A:275:VAL:HG23	1:A:309:LYS:HG2	1.98	0.46
1:A:214:PRO:HB2	1:A:215:PRO:HD3	1.96	0.46
1:A:197:ARG:O	1:A:199:ARG:HG3	2.15	0.46
1:B:265:LEU:HG	1:B:368:LEU:HB3	1.96	0.46
1:B:220:GLN:HE21	1:B:220:GLN:HA	1.80	0.46
1:A:480:GLN:NE2	1:A:610:ALA:HB3	2.30	0.46
1:A:182:THR:OG1	1:A:185:GLU:HG3	2.16	0.45
1:A:258:LEU:HD11	1:A:373:ASP:OD2	2.15	0.45
1:B:284:HIS:N	1:B:292:GLN:HE21	2.12	0.45
1:B:554:VAL:HG11	1:B:577:ALA:HB1	1.99	0.45
1:A:481:GLN:O	1:A:485:GLU:HG3	2.17	0.45
3:B:613:OCB:HAJ3	3:B:613:OCB:CAH	2.47	0.45
1:A:106:PRO:HB3	1:A:293:VAL:HG11	1.99	0.45
1:A:415:HIS:HD2	1:A:417:ASP:H	1.64	0.45
1:A:439:TYR:C	1:A:439:TYR:CD1	2.90	0.45
1:A:310:SER:O	1:A:325:CYS:HB3	2.16	0.45
1:B:129:LEU:O	1:B:133:LEU:HG	2.17	0.45
1:A:470:GLN:O	1:A:472:PRO:HD3	2.17	0.44
1:A:190:THR:O	1:A:205:VAL:HG22	2.17	0.44
1:A:490:HIS:HD2	5:A:687:HOH:O	2.01	0.44
1:A:199:ARG:NH2	1:A:367:GLU:OE1	2.39	0.44
1:B:513:ILE:O	1:B:517:GLU:HG2	2.18	0.44
1:B:36:LEU:HD21	4:B:615:MPD:HM2	1.99	0.44
1:B:123:GLU:HG3	5:B:802:HOH:O	2.16	0.44
1:B:570:ARG:CD	1:B:573:ARG:HH11	2.28	0.44
1:B:538:ASN:HA	1:B:559:VAL:HG11	1.98	0.44
1:B:453:GLU:HG3	1:B:454:THR:N	2.32	0.44
1:B:225:HIS:HE1	5:B:675:HOH:O	1.99	0.44
1:B:358:LYS:HZ3	1:B:358:LYS:HB3	1.83	0.44
1:B:338:ILE:O	1:B:342:VAL:HG23	2.18	0.44
1:A:500:GLY:HA2	5:A:617:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:PHE:CE2	1:B:367:GLU:HB2	2.52	0.44
1:A:259:ASP:OD1	1:A:261:GLU:HB3	2.18	0.44
1:B:220:GLN:NE2	1:B:220:GLN:HA	2.33	0.44
1:B:395:GLN:HG3	1:B:580:SER:O	2.18	0.43
1:B:181:LYS:HD2	1:B:187:HIS:O	2.17	0.43
1:B:561:ASN:HA	1:B:584:CYS:SG	2.57	0.43
1:B:395:GLN:HG2	1:B:582:ARG:HG2	2.00	0.43
1:B:568:HIS:CE1	1:B:570:ARG:HE	2.10	0.43
1:A:345:ARG:HA	1:A:345:ARG:NE	2.33	0.43
1:A:223:TYR:CZ	1:A:227:LYS:HD2	2.54	0.43
1:B:404:PHE:HB3	1:B:408:LEU:HD13	2.00	0.43
1:B:252:ARG:HH21	1:B:270:GLN:NE2	2.04	0.43
1:A:213:THR:OG1	1:A:215:PRO:HD2	2.18	0.43
1:A:147:HIS:H	1:A:156:MSE:SE	2.51	0.43
1:A:16:TYR:HD2	1:A:146:VAL:HG11	1.82	0.43
1:A:568:HIS:CD2	1:A:575:VAL:HB	2.54	0.43
1:B:131:HIS:HD2	5:B:742:HOH:O	2.01	0.43
1:B:553:GLY:HA3	3:B:613:OCB:HAC2	2.01	0.42
1:A:333:MSE:HG2	5:A:720:HOH:O	2.17	0.42
1:A:441:THR:OG1	2:A:613:152:H3	2.19	0.42
1:B:50:LYS:NZ	1:B:54:ILE:HD11	2.34	0.42
1:A:284:HIS:H	1:A:292:GLN:NE2	2.17	0.42
1:A:508:LEU:O	1:A:512:LEU:HG	2.19	0.42
1:A:367:GLU:HG2	1:A:369:VAL:HG22	2.01	0.42
1:A:440:GLU:HB3	1:A:455:VAL:HB	2.02	0.42
1:A:439:TYR:HB2	1:A:539:PHE:CG	2.54	0.42
1:A:570:ARG:HB2	1:A:573:ARG:HG2	2.01	0.42
1:B:103:PHE:CE1	1:B:335[A]:MSE:HE3	2.54	0.42
1:B:44:ASN:OD1	1:B:44:ASN:C	2.58	0.42
1:A:207:HIS:CG	1:A:362:ILE:HG22	2.55	0.42
1:A:604:ILE:HD13	1:A:607:MSE:CE	2.50	0.42
1:B:401:PHE:CZ	1:B:403:SER:HB2	2.55	0.42
1:A:301:ASP:OD1	1:A:303:SER:OG	2.37	0.42
1:A:268:LYS:HD2	1:A:368:LEU:CD2	2.50	0.42
1:A:297:LEU:HA	1:A:297:LEU:HD12	1.92	0.42
1:B:213:THR:HB	1:B:214:PRO:HD2	2.02	0.42
1:B:335[A]:MSE:SE	3:B:613:OCB:CAA	3.17	0.41
1:A:121:GLN:NE2	1:A:317:ALA:HA	2.35	0.41
1:A:242:SER:HB2	1:A:387:HIS:HD2	1.84	0.41
1:A:418:THR:HG23	1:A:490:HIS:HB3	2.02	0.41
1:A:326:ASP:O	1:A:329:PRO:HD2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:595:MSE:HE3	5:B:778:HOH:O	2.19	0.41
1:B:94:ARG:O	1:B:244:GLU:HG2	2.20	0.41
1:B:140:ARG:O	1:B:183:GLU:HG3	2.20	0.41
1:A:113:TYR:CE2	1:A:350:GLU:HA	2.55	0.41
1:B:230:ASN:O	1:B:231:GLU:HG3	2.21	0.41
1:B:281:SER:HB2	1:B:296:MSE:HE1	2.02	0.41
1:A:561:ASN:HA	1:A:584:CYS:SG	2.60	0.41
1:B:105:GLY:HA2	1:B:106:PRO:HD3	1.90	0.41
1:B:328:ALA:HB3	1:B:329:PRO:HD3	2.02	0.41
1:B:407:LYS:HG3	1:B:571:ASP:HB2	2.03	0.41
1:A:376:ILE:O	1:A:380:VAL:HG23	2.21	0.41
1:A:564:GLY:O	1:A:578:CYS:HA	2.20	0.41
1:B:103:PHE:HA	1:B:555:VAL:HG12	2.03	0.41
1:A:28:LEU:HD22	1:A:32:LEU:CD1	2.51	0.41
1:B:311:TYR:CD1	1:B:335[B]:MSE:HE1	2.55	0.41
1:B:213:THR:OG1	1:B:215:PRO:HD2	2.21	0.41
1:A:554:VAL:HG22	1:A:555:VAL:H	1.85	0.41
1:B:252:ARG:NH2	1:B:270:GLN:HE22	2.04	0.41
1:A:262:ASN:HD21	1:A:371:THR:HB	1.84	0.41
1:A:270:GLN:NE2	5:A:630:HOH:O	2.54	0.41
1:A:237:ILE:HD11	1:A:372:VAL:HG11	2.02	0.41
1:A:193:ALA:HA	1:A:202:VAL:HG12	2.02	0.41
1:B:408:LEU:HD23	1:B:408:LEU:C	2.41	0.41
1:A:325:CYS:SG	1:A:335:MSE:HG3	2.60	0.41
1:B:238:ALA:HB1	1:B:305:ARG:O	2.21	0.41
1:A:207:HIS:O	1:A:208:GLU:HB2	2.21	0.40
1:A:554:VAL:HG22	1:A:555:VAL:N	2.36	0.40
1:A:596:ILE:HG23	1:A:597:PHE:HD2	1.86	0.40
1:A:219:ARG:NH2	1:A:364:LEU:HD23	2.37	0.40
1:B:604:ILE:HD13	1:B:604:ILE:HA	1.99	0.40
1:A:490:HIS:HA	1:A:493:MSE:HE2	2.03	0.40
1:A:612:LEU:N	1:A:612:LEU:HD12	2.37	0.40
1:B:82:GLU:OE2	1:B:449:HIS:ND1	2.50	0.40
1:A:167:LYS:HE3	1:A:326:ASP:OD2	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	587/612 (96%)	562 (96%)	23 (4%)	2 (0%)	46	50
1	B	600/612 (98%)	581 (97%)	19 (3%)	0	100	100
All	All	1187/1224 (97%)	1143 (96%)	42 (4%)	2 (0%)	52	59

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	550	ARG
1	A	82	GLU

### 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	522/524 (100%)	510 (98%)	12 (2%)	58	71
1	B	531/524 (101%)	520 (98%)	11 (2%)	61	74
All	All	1053/1048 (100%)	1030 (98%)	23 (2%)	60	72

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	70	LEU
1	A	102	ASN
1	A	227	LYS

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Mol	Chain	Res	Type
1	A	248	TRP
1	A	261	GLU
1	A	312	ASN
1	A	313	LEU
1	A	330	TYR
1	A	439	TYR
1	A	443	MSE
1	A	504	ASP
1	B	49	LYS
1	B	102	ASN
1	B	139	LEU
1	B	248	TRP
1	B	265	LEU
1	B	297	LEU
1	B	312	ASN
1	B	313	LEU
1	B	330	TYR
1	B	434	ARG
1	B	439	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	ASN
1	A	102	ASN
1	A	109	HIS
1	A	112	HIS
1	A	178	ASN
1	A	220	GLN
1	A	270	GLN
1	A	292	GLN
1	A	378	ASN
1	A	415	HIS
1	A	470	GLN
1	A	480	GLN
1	A	490	HIS
1	A	560	HIS
1	A	568	HIS
1	A	594	GLN
1	A	608	ASN
1	B	56	GLN
1	B	87	ASN

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Mol	Chain	Res	Type
1	B	102	ASN
1	B	109	HIS
1	B	112	HIS
1	B	121	GLN
1	B	147	HIS
1	B	151	ASN
1	B	178	ASN
1	B	220	GLN
1	B	225	HIS
1	B	270	GLN
1	B	292	GLN
1	B	395	GLN
1	B	415	HIS
1	B	470	GLN
1	B	480	GLN
1	B	568	HIS
1	B	594	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

7 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	152	A	613	-	7,10,10	1.25	1 (14%)	9,14,14	1.07	1 (11%)
3	OCB	B	613	-	16,19,19	0.86	1 (6%)	20,24,24	0.97	1 (5%)
4	MPD	B	614	-	6,7,7	0.47	0	7,10,10	0.45	0
4	MPD	B	615	-	6,7,7	0.42	0	7,10,10	0.43	0
4	MPD	B	616	-	6,7,7	0.45	0	7,10,10	0.42	0
4	MPD	B	617	-	6,7,7	0.41	0	7,10,10	0.41	0
4	MPD	B	618	-	6,7,7	0.45	0	7,10,10	0.44	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
2	152	A	613	-	-	0/7/9/9	0/0/0/0
3	OCB	B	613	-	-	0/18/20/20	0/0/0/0
4	MPD	B	614	-	-	0/5/5/5	0/0/0/0
4	MPD	B	615	-	-	0/5/5/5	0/0/0/0
4	MPD	B	616	-	-	0/5/5/5	0/0/0/0
4	MPD	B	617	-	-	0/5/5/5	0/0/0/0
4	MPD	B	618	-	-	0/5/5/5	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	613	152	C4-N5	-2.42	1.47	1.51
3	B	613	OCB	CAO-NAN	-2.33	1.47	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	613	152	C3-C4-N5	-3.14	111.00	117.24
3	B	613	OCB	OAK-CAH-CAG	3.16	118.39	111.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	613	152	3	0
3	B	613	OCB	8	0
4	B	615	MPD	1	0
4	B	616	MPD	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	574/612 (93%)	0.28	39 (6%) 20 20	23, 42, 64, 91	0
1	B	583/612 (95%)	-0.18	10 (1%) 73 72	14, 26, 48, 102	0
All	All	1157/1224 (94%)	0.05	49 (4%) 40 39	14, 33, 63, 102	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	612	LEU	7.2
1	B	407	LYS	6.2
1	A	611	HIS	5.5
1	B	609	THR	4.7
1	A	610	ALA	4.5
1	A	230	ASN	4.4
1	A	601	HIS	4.4
1	B	408	LEU	4.0
1	A	233	VAL	4.0
1	B	413	ALA	3.5
1	A	609	THR	3.4
1	A	476	LEU	3.2
1	B	230	ASN	3.2
1	A	123	GLU	3.1
1	A	232	PRO	3.1
1	B	402	THR	3.1
1	A	151	ASN	3.0
1	B	404	PHE	3.0
1	A	361	ASP	2.9
1	A	228	CYS	2.7
1	A	260	PRO	2.7
1	B	610	ALA	2.7
1	A	571	ASP	2.7
1	A	372	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	229	SER	2.6
1	A	313	LEU	2.6
1	A	605	GLN	2.6
1	A	208	GLU	2.5
1	A	49	LYS	2.4
1	A	402	THR	2.4
1	A	585	PRO	2.4
1	A	358	LYS	2.4
1	A	209	GLY	2.3
1	A	234	GLY	2.3
1	A	12	ARG	2.3
1	A	275	VAL	2.3
1	A	231	GLU	2.2
1	A	116	ALA	2.2
1	A	202	VAL	2.2
1	A	371	THR	2.2
1	B	406	LYS	2.1
1	A	604	ILE	2.1
1	A	608	ASN	2.1
1	A	312	ASN	2.1
1	A	117	ARG	2.1
1	A	187	HIS	2.1
1	B	409	THR	2.0
1	A	311	TYR	2.0
1	A	286	THR	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates 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. 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
4	MPD	B	615	8/8	0.83	0.33	13.78	68,68,72,72	0
4	MPD	B	616	8/8	0.74	0.33	9.19	68,69,70,70	0
4	MPD	B	614	8/8	0.90	0.21	4.14	48,50,52,53	0
4	MPD	B	618	8/8	0.78	0.31	2.75	75,76,77,77	0
2	152	A	613	11/11	0.86	0.19	2.19	28,33,35,36	0
3	OCB	B	613	20/20	0.95	0.17	0.87	20,24,36,37	0
4	MPD	B	617	8/8	0.70	0.34	-	62,64,65,66	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.