Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 01:13 PM EDT

PDB ID	:	1XL7
Title	:	Crystal Structure of Mouse Carnitine Octanoyltransferase
Authors	:	Jogl, G.; Hsiao, Y.S.; Tong, L.
Deposited on	:	2004-09-30
Resolution	:	2.00 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 (i) 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 (i)) were used in the production of this report:

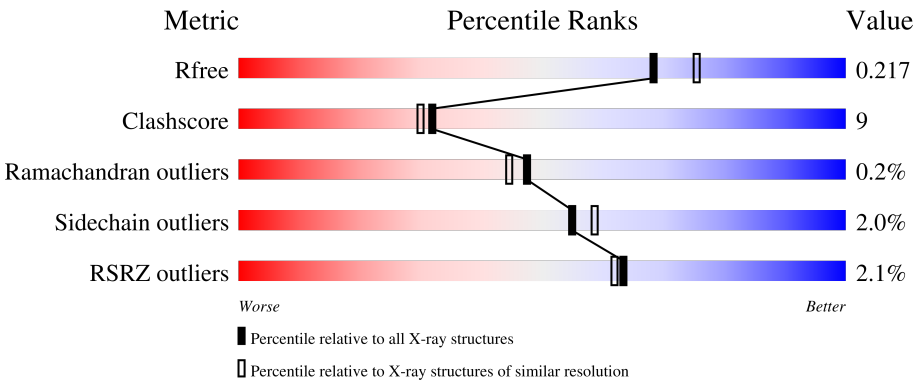
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 <=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	% 77% 19% •
1	B	612	3% 82% 15% • •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 10260 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
			Total	C	N	O	S	Se			
1	A	591	4776	3042	829	872	16	17	0	0	0
1	B	600	4843	3085	839	886	16	17	0	0	0

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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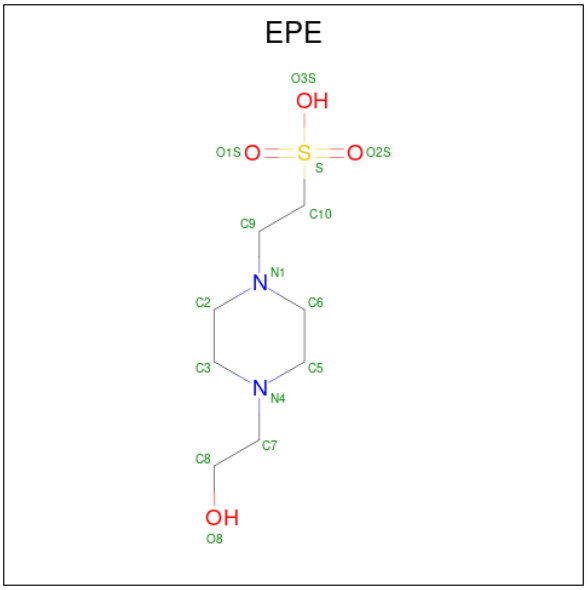
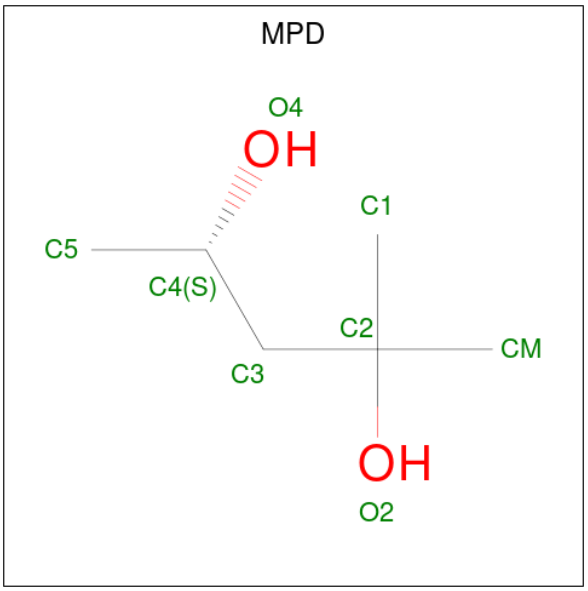
Chain	Residue	Modelled	Actual	Comment	Reference
B	443	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	469	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	483	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	493	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	494	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	558	MSE	MET	MODIFIED RESIDUE	UNP Q9DC50
B	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 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	15	8	2	4	1	0	0
			Total	C	N	O	S		
2	B	1	15	8	2	4	1	0	0

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



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

- Molecule 4 is water.

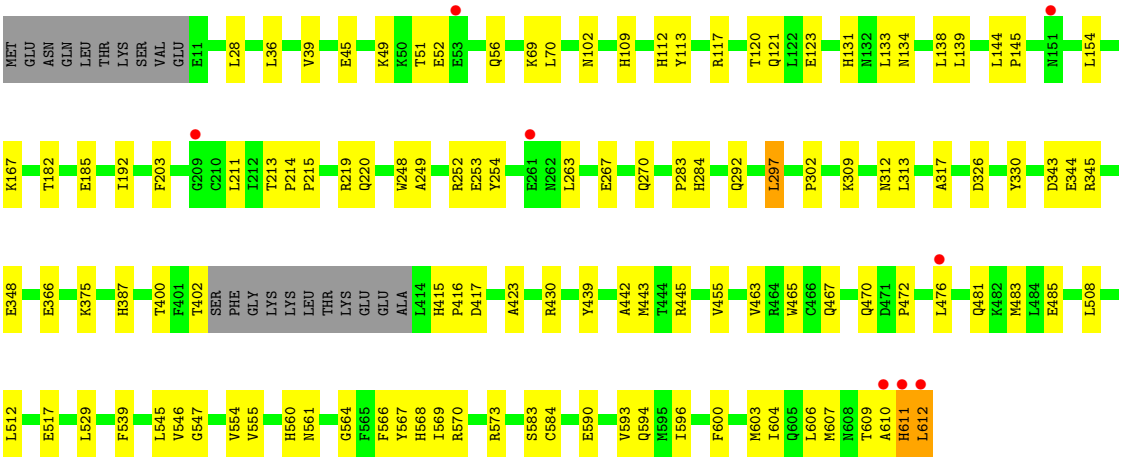
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	229	229	229	0	0
4	B	310	310	310	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

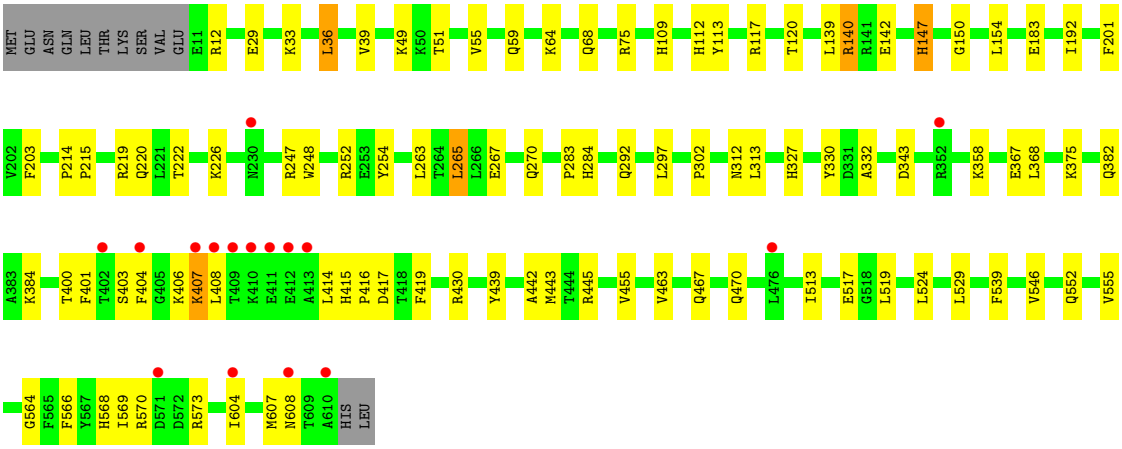
• Molecule 1: Peroxisomal carnitine O-octanoyltransferase

Chain A:



• Molecule 1: Peroxisomal carnitine O-octanoyltransferase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	162.63Å 162.63Å 158.58Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.92 – 2.00 29.92 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.6 (29.92-2.00) 98.3 (29.92-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$< I/\sigma(I) >$ [1]	2.38 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.190 , 0.217 0.192 , 0.217	Depositor DCC
R_{free} test set	14604 reflections (6.93%)	wwPDB-VP
Wilson B-factor (Å^2)	20.9	Xtriage
Anisotropy	0.065	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 56.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.016 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	10260	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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

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	0.32	0/4884	0.57	1/6585 (0.0%)
1	B	0.35	0/4952	0.59	2/6676 (0.0%)
All	All	0.33	0/9836	0.58	3/13261 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	564	GLY	N-CA-C	-6.04	98.00	113.10
1	A	564	GLY	N-CA-C	-5.76	98.71	113.10
1	B	150	GLY	N-CA-C	-5.63	99.04	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	85	0
1	B	4843	0	4750	84	0
2	A	15	0	18	0	0
2	B	15	0	18	1	0
3	A	16	0	28	0	0
3	B	56	0	98	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	229	0	0	6	0
4	B	310	0	0	5	0
All	All	10260	0	9588	171	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 9.

All (171) 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:283:PRO:HA	1:A:292:GLN:HE21	1.28	0.93
2:B:701:EPE:H101	3:B:702:MPD:H51	1.49	0.93
1:B:604:ILE:HA	1:B:607:MSE:HE3	1.54	0.89
1:A:252:ARG:HH21	1:A:270:GLN:HE22	1.21	0.86
1:A:604:ILE:HA	1:A:607:MSE:HE3	1.58	0.85
1:B:252:ARG:HH21	1:B:270:GLN:HE22	1.24	0.85
1:A:416:PRO:HB2	1:A:569:ILE:HD13	1.64	0.78
1:B:406:LYS:HB3	1:B:569:ILE:HG21	1.67	0.77
1:A:568:HIS:HE1	1:A:570:ARG:HE	1.32	0.77
1:B:219:ARG:HH21	1:B:220:GLN:HE22	1.31	0.75
1:A:134:ASN:O	1:A:138:LEU:HD13	1.86	0.74
1:B:430:ARG:HE	1:B:470:GLN:HE22	1.36	0.74
1:A:52:GLU:O	1:A:56:GLN:HG3	1.89	0.73
1:B:112:HIS:CD2	1:B:113:TYR:H	2.07	0.73
1:B:404:PHE:CE2	1:B:604:ILE:HG13	2.25	0.72
1:A:465:TRP:CZ2	1:A:483:MSE:HE2	2.26	0.71
1:B:283:PRO:HA	1:B:292:GLN:HE21	1.56	0.71
1:A:430:ARG:HE	1:A:470:GLN:NE2	1.90	0.70
1:B:415:HIS:HD2	1:B:417:ASP:H	1.40	0.70
1:B:430:ARG:HE	1:B:470:GLN:NE2	1.89	0.69
1:B:568:HIS:HE1	1:B:570:ARG:HE	1.39	0.69
1:A:570:ARG:HD2	1:A:573:ARG:HH11	1.58	0.69
1:B:284:HIS:H	1:B:292:GLN:NE2	1.92	0.67
1:A:123:GLU:CD	1:A:123:GLU:H	1.97	0.66
1:A:430:ARG:HE	1:A:470:GLN:HE22	1.42	0.65
1:A:609:THR:O	1:A:612:LEU:HD13	1.97	0.65
1:A:465:TRP:HZ2	1:A:483:MSE:HE2	1.62	0.65
1:A:546:VAL:HG23	1:A:568:HIS:HB3	1.76	0.65
1:A:283:PRO:HA	1:A:292:GLN:NE2	2.06	0.63
1:A:252:ARG:NH2	1:A:270:GLN:HE22	1.95	0.62
1:A:463:VAL:O	1:A:467:GLN:HG3	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:406:LYS:HB3	1:B:569:ILE:CG2	2.30	0.61
1:A:252:ARG:HH21	1:A:270:GLN:NE2	1.94	0.61
1:B:604:ILE:HD13	1:B:607:MSE:CE	2.31	0.60
1:A:604:ILE:HD13	1:A:607:MSE:CE	2.31	0.60
1:A:117:ARG:HE	1:A:120:THR:HG21	1.67	0.59
1:A:284:HIS:H	1:A:292:GLN:NE2	1.99	0.59
1:B:302:PRO:HG2	1:B:384:LYS:HG3	1.84	0.59
1:A:219:ARG:HH21	1:A:220:GLN:HE22	1.51	0.58
1:A:483:MSE:HE3	1:A:606:LEU:HD23	1.84	0.58
1:B:604:ILE:HD13	1:B:607:MSE:HE3	1.85	0.58
1:A:219:ARG:HH21	1:A:220:GLN:NE2	2.01	0.58
1:B:401:PHE:CE2	1:B:403:SER:HB2	2.39	0.57
1:A:545:LEU:HG	1:A:547:GLY:H	1.70	0.57
1:A:546:VAL:CG2	1:A:568:HIS:HB3	2.35	0.57
1:B:222:THR:O	1:B:226:LYS:HG2	2.04	0.57
1:A:600:PHE:O	1:A:604:ILE:HG12	2.05	0.57
1:A:112:HIS:CD2	1:A:113:TYR:H	2.22	0.57
1:A:604:ILE:HA	1:A:607:MSE:CE	2.32	0.56
1:A:604:ILE:HD13	1:A:607:MSE:HE1	1.88	0.56
1:B:33:LYS:HE2	3:B:705:MPD:H31	1.87	0.56
1:B:219:ARG:HH21	1:B:220:GLN:NE2	2.02	0.56
1:B:332:ALA:HB2	3:B:702:MPD:H13	1.87	0.56
1:B:154:LEU:HD13	1:B:443:MSE:HE1	1.86	0.56
1:B:59:GLN:NE2	3:B:705:MPD:H51	2.21	0.55
1:B:192:ILE:HD11	1:B:203:PHE:CZ	2.41	0.55
1:B:415:HIS:CD2	1:B:417:ASP:H	2.21	0.55
1:A:249:ALA:O	1:A:253:GLU:HG3	2.06	0.55
1:B:406:LYS:HG3	1:B:407:LYS:N	2.22	0.55
1:A:192:ILE:HD11	1:A:203:PHE:CZ	2.41	0.55
1:B:147:HIS:CD2	1:B:445:ARG:HH22	2.25	0.55
1:B:400:THR:CG2	1:B:573:ARG:HD3	2.37	0.54
1:A:443:MSE:HE2	1:A:445:ARG:HD3	1.89	0.54
1:B:524:LEU:C	1:B:524:LEU:HD23	2.28	0.54
1:B:404:PHE:CZ	1:B:604:ILE:HG13	2.42	0.54
1:A:167:LYS:HE3	1:A:326:ASP:OD2	2.08	0.54
1:B:546:VAL:HG12	1:B:546:VAL:O	2.06	0.54
1:B:112:HIS:HD2	1:B:113:TYR:H	1.56	0.53
1:B:254:TYR:CZ	1:B:375:LYS:HD3	2.44	0.53
1:B:265:LEU:HG	1:B:368:LEU:HB3	1.91	0.52
1:B:513:ILE:O	1:B:517:GLU:HG2	2.10	0.52
1:B:416:PRO:HB2	1:B:569:ILE:HD13	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:LEU:O	1:A:267:GLU:HG3	2.09	0.52
1:B:358:LYS:NZ	1:B:358:LYS:HB3	2.25	0.52
1:A:345:ARG:HH21	1:A:348:GLU:CD	2.13	0.51
1:B:404:PHE:HB2	1:B:408:LEU:HD22	1.92	0.51
1:A:546:VAL:HG22	1:A:567:TYR:O	2.11	0.51
1:B:252:ARG:HH21	1:B:270:GLN:NE2	2.02	0.51
1:A:214:PRO:HB2	1:A:215:PRO:HD3	1.92	0.51
1:B:414:LEU:HD12	1:B:607:MSE:SE	2.60	0.51
1:A:483:MSE:CE	1:A:606:LEU:HD23	2.39	0.51
1:A:213:THR:OG1	1:A:215:PRO:HD2	2.11	0.50
1:B:36:LEU:CD2	1:B:55:VAL:HG11	2.40	0.50
1:B:517:GLU:HB2	1:B:519:LEU:HG	1.93	0.50
1:A:442:ALA:HB2	1:A:455:VAL:HG23	1.94	0.50
1:B:404:PHE:CB	1:B:408:LEU:HD22	2.42	0.50
1:B:463:VAL:O	1:B:467:GLN:HG3	2.11	0.50
1:B:117:ARG:HE	1:B:120:THR:HG21	1.77	0.49
1:B:419:PHE:HE1	1:B:607:MSE:HE1	1.77	0.49
1:B:263:LEU:O	1:B:267:GLU:HG3	2.13	0.49
1:B:75:ARG:HA	3:B:706:MPD:HM1	1.94	0.49
1:B:401:PHE:CZ	1:B:403:SER:HB2	2.47	0.49
1:B:442:ALA:HB2	1:B:455:VAL:HG23	1.95	0.49
1:A:121:GLN:NE2	1:A:317:ALA:HA	2.27	0.49
1:A:560:HIS:HE1	1:A:583:SER:OG	1.95	0.49
1:B:29:GLU:O	1:B:33:LYS:HG2	2.13	0.49
1:B:75:ARG:HG2	3:B:706:MPD:HM1	1.94	0.48
1:B:201:PHE:CE2	1:B:367:GLU:HB2	2.48	0.48
1:B:252:ARG:NH2	1:B:270:GLN:HE22	2.04	0.48
1:A:400:THR:CG2	1:A:573:ARG:HD3	2.42	0.48
1:A:517:GLU:HG3	4:A:825:HOH:O	2.13	0.48
1:A:254:TYR:CE2	1:A:375:LYS:HD3	2.49	0.47
1:B:214:PRO:HB2	1:B:215:PRO:HD3	1.95	0.47
1:A:402:THR:HG22	1:A:573:ARG:HB3	1.95	0.47
1:A:109:HIS:HD2	1:A:343:ASP:OD1	1.96	0.47
1:B:36:LEU:HD22	1:B:55:VAL:HG11	1.97	0.47
1:B:49:LYS:HE2	4:B:933:HOH:O	2.15	0.47
1:A:555:VAL:HG22	1:A:566:PHE:CZ	2.50	0.47
1:B:401:PHE:HE2	1:B:404:PHE:CZ	2.33	0.47
1:A:590:GLU:O	1:A:594:GLN:HG3	2.15	0.46
1:A:133:LEU:HD13	1:A:211:LEU:HD11	1.97	0.46
1:A:344:GLU:O	1:A:348:GLU:HG3	2.16	0.46
1:B:140:ARG:O	1:B:183:GLU:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:406:LYS:C	1:B:408:LEU:H	2.19	0.46
1:A:481:GLN:O	1:A:485:GLU:HG3	2.15	0.46
1:B:64:LYS:O	1:B:68:GLN:HG3	2.15	0.46
1:B:400:THR:HG21	1:B:573:ARG:HD3	1.97	0.46
1:A:416:PRO:CB	1:A:569:ILE:HD13	2.42	0.46
1:B:36:LEU:O	1:B:39:VAL:HG22	2.16	0.46
1:A:302:PRO:HG3	1:A:387:HIS:ND1	2.31	0.45
1:A:443:MSE:CE	1:A:445:ARG:HB2	2.46	0.45
1:A:182:THR:OG1	1:A:185:GLU:HG3	2.17	0.45
1:A:51:THR:OG1	1:A:529:LEU:HD12	2.17	0.45
1:A:45:GLU:O	1:A:49:LYS:HG2	2.16	0.45
1:B:247:ARG:NH2	1:B:382:GLN:NE2	2.65	0.45
3:B:702:MPD:H53	4:B:1018:HOH:O	2.17	0.45
1:A:154:LEU:HD13	1:A:443:MSE:HE1	1.99	0.44
1:B:568:HIS:CE1	1:B:570:ARG:HE	2.28	0.44
1:A:476:LEU:HD21	1:A:609:THR:HG22	2.00	0.44
1:B:524:LEU:HD23	1:B:524:LEU:O	2.16	0.44
1:A:131:HIS:HD2	4:A:770:HOH:O	2.01	0.44
1:A:554:VAL:HG23	1:A:566:PHE:HD2	1.83	0.44
1:A:470:GLN:O	1:A:472:PRO:HD3	2.17	0.44
1:B:12:ARG:NH2	1:B:142:GLU:OE2	2.51	0.44
1:A:593:VAL:O	1:A:596:ILE:HG22	2.18	0.43
1:B:439:TYR:HB2	1:B:539:PHE:CG	2.53	0.43
1:A:366:GLU:HG3	4:A:738:HOH:O	2.17	0.43
1:B:415:HIS:HD2	1:B:417:ASP:N	2.09	0.43
1:B:109:HIS:HD2	1:B:343:ASP:OD1	2.01	0.43
1:A:611:HIS:O	1:A:612:LEU:OXT	2.36	0.43
1:A:309:LYS:HE3	4:A:622:HOH:O	2.19	0.43
1:A:144:LEU:HD12	1:A:145:PRO:HD2	2.01	0.43
1:A:387:HIS:HD2	4:A:659:HOH:O	2.02	0.43
1:B:604:ILE:HD13	1:B:607:MSE:HE1	2.00	0.43
1:A:36:LEU:O	1:A:39:VAL:HG22	2.19	0.42
1:B:51:THR:OG1	1:B:529:LEU:HD12	2.19	0.42
1:B:439:TYR:CD1	1:B:439:TYR:C	2.93	0.42
1:A:561:ASN:HA	1:A:584:CYS:SG	2.59	0.42
1:A:439:TYR:CD1	1:A:439:TYR:C	2.93	0.42
1:A:254:TYR:CZ	1:A:375:LYS:HD3	2.55	0.42
1:B:406:LYS:HB2	1:B:416:PRO:HG3	2.02	0.42
1:A:423:ALA:HB2	1:A:603:MSE:SE	2.69	0.42
1:A:220:GLN:NE2	1:A:220:GLN:HA	2.34	0.42
1:A:439:TYR:HB2	1:A:539:PHE:CG	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:430:ARG:NE	1:B:470:GLN:HE22	2.11	0.42
1:B:552:GLN:OE1	1:B:568:HIS:HD2	2.03	0.42
1:A:603:MSE:O	1:A:607:MSE:HG3	2.19	0.41
1:B:419:PHE:CE1	1:B:607:MSE:HE1	2.54	0.41
1:B:414:LEU:CD1	1:B:607:MSE:SE	3.18	0.41
1:B:555:VAL:HG22	1:B:566:PHE:CZ	2.55	0.41
1:B:117:ARG:NH2	4:B:917:HOH:O	2.53	0.41
1:A:508:LEU:O	1:A:512:LEU:HG	2.20	0.41
1:B:327:HIS:NE2	3:B:702:MPD:H52	2.36	0.41
1:A:568:HIS:CE1	1:A:570:ARG:HE	2.24	0.41
1:B:147:HIS:HE1	4:B:869:HOH:O	2.02	0.41
1:B:604:ILE:HG22	1:B:608:ASN:ND2	2.36	0.41
1:A:131:HIS:HE1	4:A:732:HOH:O	2.04	0.41
1:A:297:LEU:HD12	1:A:297:LEU:HA	1.90	0.41
1:A:415:HIS:CD2	1:A:417:ASP:H	2.40	0.40
1:B:247:ARG:NH2	4:B:1013:HOH:O	2.55	0.40
1:A:610:ALA:O	1:A:611:HIS:C	2.60	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	587/612 (96%)	569 (97%)	17 (3%)	1 (0%)	47	44
1	B	598/612 (98%)	583 (98%)	14 (2%)	1 (0%)	47	44
All	All	1185/1224 (97%)	1152 (97%)	31 (3%)	2 (0%)	47	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	407	LYS

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Mol	Chain	Res	Type
1	A	611	HIS

5.3.2 Protein sidechains (i)

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%)	511 (98%)	11 (2%)	53	57
1	B	529/524 (101%)	519 (98%)	10 (2%)	57	61
All	All	1051/1048 (100%)	1030 (98%)	21 (2%)	55	58

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	69	LYS
1	A	70	LEU
1	A	102	ASN
1	A	139	LEU
1	A	248	TRP
1	A	297	LEU
1	A	312	ASN
1	A	313	LEU
1	A	330	TYR
1	A	612	LEU
1	B	36	LEU
1	B	139	LEU
1	B	140	ARG
1	B	147	HIS
1	B	248	TRP
1	B	265	LEU
1	B	297	LEU
1	B	312	ASN
1	B	313	LEU
1	B	330	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	A	87	ASN
1	A	102	ASN
1	A	109	HIS
1	A	112	HIS
1	A	121	GLN
1	A	131	HIS
1	A	157	ASN
1	A	178	ASN
1	A	220	GLN
1	A	225	HIS
1	A	270	GLN
1	A	292	GLN
1	A	378	ASN
1	A	382	GLN
1	A	387	HIS
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	B	56	GLN
1	B	87	ASN
1	B	102	ASN
1	B	109	HIS
1	B	112	HIS
1	B	121	GLN
1	B	131	HIS
1	B	147	HIS
1	B	151	ASN
1	B	178	ASN
1	B	220	GLN
1	B	270	GLN
1	B	292	GLN
1	B	382	GLN
1	B	395	GLN
1	B	415	HIS
1	B	470	GLN
1	B	490	HIS

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Mol	Chain	Res	Type
1	B	568	HIS
1	B	594	GLN
1	B	608	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

11 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	EPE	A	613	-	15,15,15	1.36	1 (6%)	19,20,20	0.59	0
3	MPD	B	707	-	7,7,7	0.45	0	9,10,10	0.43	0
3	MPD	A	615	-	7,7,7	0.45	0	9,10,10	0.41	0
3	MPD	B	706	-	7,7,7	0.52	0	9,10,10	0.50	0
3	MPD	B	703	-	7,7,7	0.46	0	9,10,10	0.51	0
3	MPD	B	702	-	7,7,7	0.45	0	9,10,10	0.60	0
3	MPD	B	704	-	7,7,7	0.49	0	9,10,10	0.44	0
3	MPD	B	705	-	7,7,7	0.43	0	9,10,10	0.44	0
3	MPD	A	614	-	7,7,7	0.46	0	9,10,10	0.49	0
3	MPD	B	708	-	7,7,7	0.50	0	9,10,10	0.45	0
2	EPE	B	701	-	15,15,15	1.35	1 (6%)	19,20,20	0.67	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	EPE	A	613	-	-	0/9/19/19	0/1/1/1
3	MPD	B	707	-	-	3/5/5/5	-
3	MPD	A	615	-	-	2/5/5/5	-
3	MPD	B	706	-	-	0/5/5/5	-
3	MPD	B	703	-	-	0/5/5/5	-
3	MPD	B	702	-	-	2/5/5/5	-
3	MPD	B	704	-	-	3/5/5/5	-
3	MPD	B	705	-	-	5/5/5/5	-
3	MPD	A	614	-	-	0/5/5/5	-
3	MPD	B	708	-	-	2/5/5/5	-
2	EPE	B	701	-	-	0/9/19/19	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	613	EPE	C10-S	4.25	1.83	1.77
2	B	701	EPE	C10-S	4.15	1.83	1.77

There are no bond angle outliers.

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	615	MPD	C2-C3-C4-O4
3	A	615	MPD	C2-C3-C4-C5
3	B	702	MPD	C2-C3-C4-O4
3	B	702	MPD	C2-C3-C4-C5
3	B	705	MPD	C2-C3-C4-O4
3	B	705	MPD	C2-C3-C4-C5
3	B	704	MPD	C1-C2-C3-C4
3	B	704	MPD	CM-C2-C3-C4
3	B	705	MPD	C1-C2-C3-C4
3	B	707	MPD	C1-C2-C3-C4
3	B	707	MPD	CM-C2-C3-C4
3	B	708	MPD	C1-C2-C3-C4
3	B	704	MPD	O2-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
3	B	705	MPD	O2-C2-C3-C4
3	B	707	MPD	O2-C2-C3-C4
3	B	705	MPD	CM-C2-C3-C4
3	B	708	MPD	CM-C2-C3-C4

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	706	MPD	2	0
3	B	702	MPD	4	0
3	B	705	MPD	2	0
2	B	701	EPE	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	574/612 (93%)	-0.18	8 (1%)	75	74	15, 28, 45, 70	0
1	B	583/612 (95%)	-0.36	16 (2%)	54	53	11, 19, 40, 99	0
All	All	1157/1224 (94%)	-0.27	24 (2%)	63	62	11, 24, 43, 99	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	408	LEU	8.4
1	B	407	LYS	7.8
1	A	612	LEU	5.6
1	B	404	PHE	5.1
1	B	402	THR	4.8
1	B	413	ALA	4.7
1	A	611	HIS	4.5
1	B	411	GLU	3.0
1	B	409	THR	3.0
1	A	151	ASN	2.9
1	B	412	GLU	2.9
1	A	610	ALA	2.9
1	A	476	LEU	2.8
1	A	209	GLY	2.6
1	B	230	ASN	2.6
1	B	410	LYS	2.6
1	B	610	ALA	2.5
1	A	53	GLU	2.4
1	B	476	LEU	2.1
1	B	604	ILE	2.1
1	B	608	ASN	2.1
1	B	352	ARG	2.1
1	B	571	ASP	2.0
1	A	261	GLU	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 monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	MPD	B	708	8/8	0.67	0.23	64,65,65,66	0
3	MPD	B	707	8/8	0.80	0.17	49,50,50,51	0
3	MPD	A	615	8/8	0.81	0.17	28,31,33,35	0
3	MPD	B	704	8/8	0.81	0.17	61,61,62,63	0
3	MPD	B	706	8/8	0.83	0.21	52,52,54,54	0
3	MPD	B	705	8/8	0.85	0.16	40,42,44,45	0
3	MPD	B	702	8/8	0.85	0.16	18,23,26,30	0
3	MPD	A	614	8/8	0.89	0.22	34,36,39,40	0
3	MPD	B	703	8/8	0.92	0.13	38,39,40,41	0
2	EPE	B	701	15/15	0.96	0.14	17,24,38,42	0
2	EPE	A	613	15/15	0.96	0.11	23,29,39,42	0

6.5 Other polymers (i)

There are no such residues in this entry.