



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 08:25 am BST

PDB ID : 2ZYR
Title : A. Fulgidus lipase with fatty acid fragment and magnesium
Authors : Chen, C.K.; Ko, T.P.; Guo, R.T.; Wang, A.H.
Deposited on : 2009-01-28
Resolution : 1.77 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

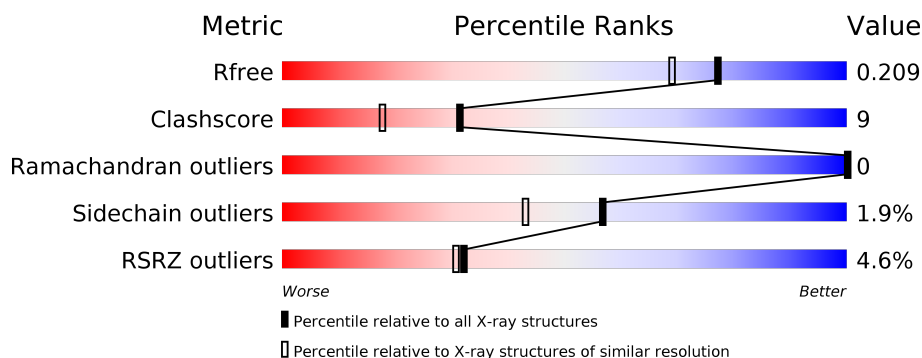
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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 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	484	<div> <div>4%</div> <div> <div></div> <div>81%</div> <div>14%</div> <div>• •</div> </div> </div>
1	B	484	<div> <div>5%</div> <div> <div></div> <div>77%</div> <div>15%</div> <div>• 6%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8535 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.

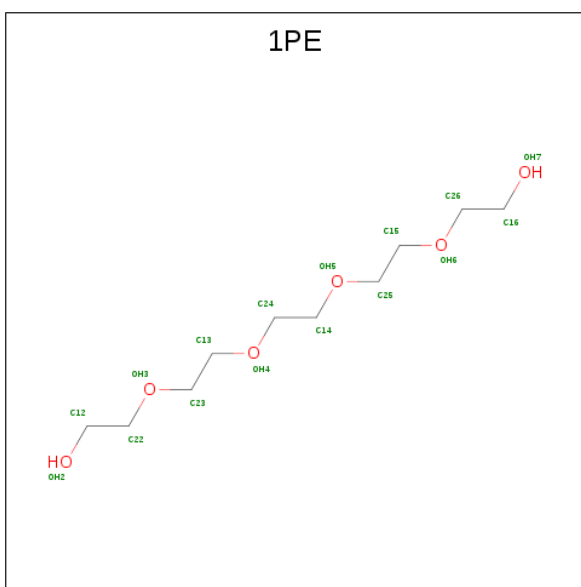
- Molecule 1 is a protein called Lipase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3660	2357	606	685	12			
1	B	454	Total	C	N	O	S	0	0	0
			3586	2310	593	671	12			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	475	VAL	-	EXPRESSION TAG	UNP O28511
A	476	ASP	-	EXPRESSION TAG	UNP O28511
A	477	LYS	-	EXPRESSION TAG	UNP O28511
A	478	LEU	-	EXPRESSION TAG	UNP O28511
A	479	ALA	-	EXPRESSION TAG	UNP O28511
A	480	ALA	-	EXPRESSION TAG	UNP O28511
A	481	ALA	-	EXPRESSION TAG	UNP O28511
A	482	LEU	-	EXPRESSION TAG	UNP O28511
A	483	GLU	-	EXPRESSION TAG	UNP O28511
A	484	HIS	-	EXPRESSION TAG	UNP O28511
B	475	VAL	-	EXPRESSION TAG	UNP O28511
B	476	ASP	-	EXPRESSION TAG	UNP O28511
B	477	LYS	-	EXPRESSION TAG	UNP O28511
B	478	LEU	-	EXPRESSION TAG	UNP O28511
B	479	ALA	-	EXPRESSION TAG	UNP O28511
B	480	ALA	-	EXPRESSION TAG	UNP O28511
B	481	ALA	-	EXPRESSION TAG	UNP O28511
B	482	LEU	-	EXPRESSION TAG	UNP O28511
B	483	GLU	-	EXPRESSION TAG	UNP O28511
B	484	HIS	-	EXPRESSION TAG	UNP O28511

- Molecule 2 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			16	10	6		
2	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Mg	0	0
			2	2		
3	A	3	Total	Mg	0	0
			3	3		

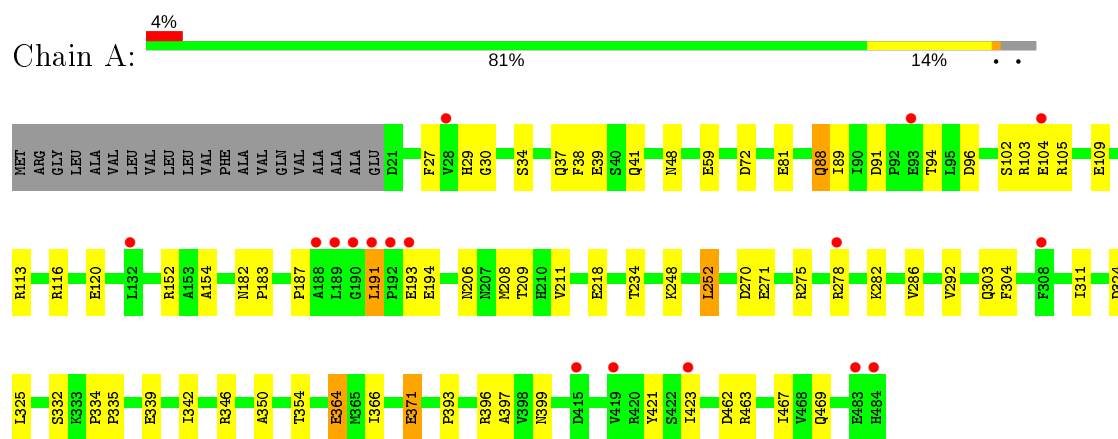
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	697	Total	O	0	0
			697	697		
4	B	558	Total	O	0	0
			558	558		

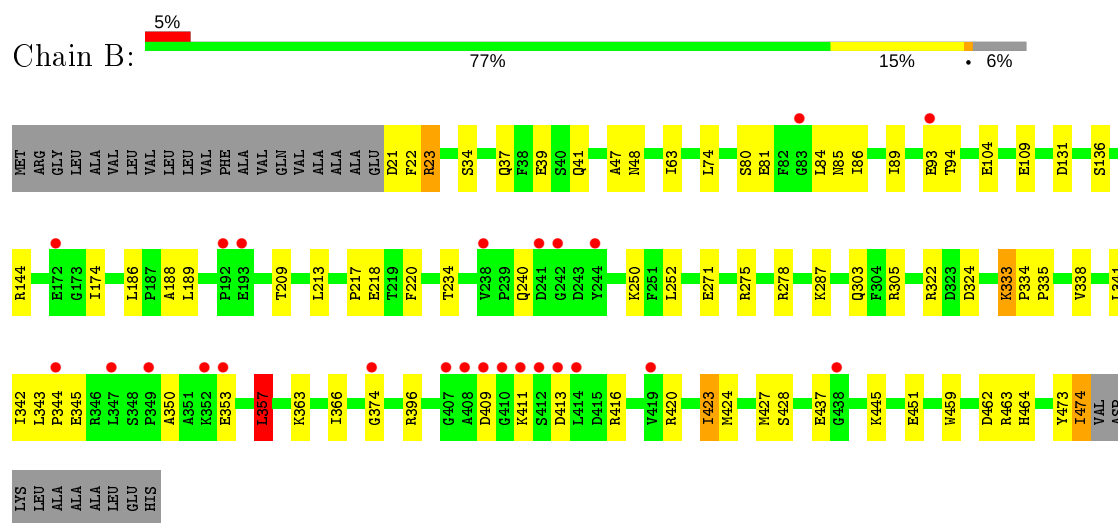
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Lipase, putative



- Molecule 1: Lipase, putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.29Å 106.67Å 175.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.77 29.40 – 1.77	Depositor EDS
% Data completeness (in resolution range)	93.1 (30.00-1.77) 93.7 (29.40-1.77)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 1.77Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.178 , 0.214 0.173 , 0.209	Depositor DCC
R_{free} test set	4527 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	0.157	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8535	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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: MG, 1PE

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	0.95	2/3748 (0.1%)	0.93	5/5088 (0.1%)
1	B	0.85	0/3673	0.89	5/4986 (0.1%)
All	All	0.90	2/7421 (0.0%)	0.91	10/10074 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	154	ALA	CA-CB	5.56	1.64	1.52
1	A	364	GLU	CG-CD	5.27	1.59	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	213	LEU	CA-CB-CG	6.31	129.81	115.30
1	B	23	ARG	NE-CZ-NH2	-6.11	117.24	120.30
1	A	113	ARG	NE-CZ-NH2	-5.91	117.35	120.30
1	B	374	GLY	N-CA-C	5.58	127.04	113.10
1	B	363	LYS	N-CA-C	-5.31	96.67	111.00
1	A	462	ASP	CB-CG-OD1	5.27	123.04	118.30
1	A	270	ASP	CB-CG-OD1	5.22	123.00	118.30
1	B	357	LEU	CA-CB-CG	5.09	127.01	115.30
1	A	334	PRO	N-CA-C	-5.09	98.87	112.10
1	A	152	ARG	NE-CZ-NH2	-5.05	117.78	120.30

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	3660	0	3615	65	0
1	B	3586	0	3539	66	0
2	A	16	0	22	0	0
2	B	13	0	17	0	0
3	A	3	0	0	0	0
3	B	2	0	0	0	0
4	A	697	0	0	17	0
4	B	558	0	0	21	0
All	All	8535	0	7193	130	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 (130) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:287:LYS:HE2	4:B:1031:HOH:O	1.55	1.05
1:A:105:ARG:O	1:A:109:GLU:HG3	1.64	0.97
1:B:424:MET:HB3	1:B:427:MET:HE3	1.54	0.89
1:A:193:GLU:HG3	4:A:1087:HOH:O	1.77	0.85
1:B:423:ILE:HD11	4:B:1123:HOH:O	1.82	0.79
1:A:120:GLU:HG3	4:A:1054:HOH:O	1.84	0.77
1:B:85:ASN:HB2	4:B:1121:HOH:O	1.83	0.77
1:B:287:LYS:HD3	4:B:1027:HOH:O	1.84	0.77
1:B:85:ASN:HB3	4:B:499:HOH:O	1.84	0.77
1:B:420:ARG:HD3	4:B:672:HOH:O	1.84	0.77
1:B:445:LYS:HG2	1:B:451:GLU:HG2	1.67	0.76
1:A:194:GLU:HA	4:A:738:HOH:O	1.85	0.76
1:A:463:ARG:HD3	4:A:790:HOH:O	1.90	0.72
1:A:282:LYS:HG3	1:A:292:VAL:HG21	1.73	0.71
1:A:354:THR:HG21	4:A:811:HOH:O	1.91	0.70
1:A:423:ILE:HG23	4:A:1016:HOH:O	1.89	0.70
1:A:423:ILE:CG2	4:A:1016:HOH:O	2.40	0.69
1:A:103:ARG:HG2	1:A:103:ARG:HH11	1.56	0.69
1:B:39:GLU:OE2	4:B:513:HOH:O	2.10	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:GLU:HG2	4:B:641:HOH:O	1.92	0.69
1:A:72:ASP:OD2	1:A:103:ARG:NH2	2.26	0.68
1:A:103:ARG:NH1	1:A:103:ARG:HG2	2.08	0.68
1:B:462:ASP:OD2	1:B:463:ARG:HG3	1.95	0.67
1:A:34:SER:H	1:A:37:GLN:HE22	1.43	0.67
1:A:89:ILE:HG22	1:A:89:ILE:O	1.94	0.66
1:A:271:GLU:H	1:A:271:GLU:CD	1.96	0.66
1:B:186:LEU:HD21	1:B:189:LEU:HD12	1.75	0.66
1:B:437:GLU:O	4:B:947:HOH:O	2.14	0.66
1:A:88:GLN:HG2	1:A:89:ILE:HD13	1.78	0.65
1:A:120:GLU:CG	4:A:1054:HOH:O	2.43	0.65
1:A:102:SER:HB3	1:A:105:ARG:HB3	1.79	0.64
1:B:424:MET:HB3	1:B:427:MET:CE	2.27	0.64
1:B:240:GLN:HG2	1:B:322:ARG:HA	1.80	0.64
1:A:286:VAL:HG21	1:A:304:PHE:HE1	1.62	0.64
1:B:209:THR:HG21	1:B:366:ILE:HD11	1.80	0.64
1:B:34:SER:H	1:B:37:GLN:HE22	1.46	0.62
1:B:48:ASN:HD21	1:B:234:THR:H	1.47	0.61
1:B:338:VAL:HG12	1:B:424:MET:CE	2.31	0.60
1:A:278:ARG:HD3	4:A:779:HOH:O	2.01	0.60
1:A:350:ALA:O	1:A:354:THR:HG23	2.02	0.59
1:A:282:LYS:HG3	1:A:292:VAL:CG2	2.32	0.58
1:B:21:ASP:N	4:B:705:HOH:O	2.36	0.58
1:B:93:GLU:HG3	4:B:787:HOH:O	2.04	0.57
1:A:371:GLU:O	1:A:371:GLU:CD	2.44	0.56
1:B:275:ARG:HH22	1:B:303:GLN:NE2	2.03	0.56
1:A:423:ILE:HG12	1:A:423:ILE:O	2.06	0.56
1:B:21:ASP:CG	1:B:22:PHE:H	2.09	0.56
1:B:81:GLU:HG2	1:B:188:ALA:HB1	1.88	0.56
1:B:74:LEU:HD23	1:B:144:ARG:HG2	1.89	0.55
1:B:473:TYR:OH	4:B:1202:HOH:O	2.18	0.54
1:A:206:ASN:ND2	4:A:733:HOH:O	2.40	0.53
1:B:80:SER:O	1:B:84:LEU:HG	2.07	0.53
1:A:286:VAL:CG2	1:A:304:PHE:HE1	2.21	0.53
1:B:357:LEU:N	1:B:357:LEU:HD23	2.23	0.53
1:A:396:ARG:NH2	4:A:1152:HOH:O	2.41	0.53
1:B:338:VAL:HG12	1:B:424:MET:HE1	1.90	0.52
1:A:467:ILE:N	1:A:467:ILE:HD12	2.25	0.52
1:A:81:GLU:HB3	4:A:979:HOH:O	2.09	0.52
1:A:208:MET:HE1	1:A:218:GLU:HB2	1.91	0.51
1:B:186:LEU:CD2	1:B:189:LEU:HD12	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:HIS:HD2	1:A:30:GLY:O	1.94	0.51
1:B:333:LYS:HG3	1:B:334:PRO:HD2	1.93	0.50
1:B:462:ASP:CG	1:B:463:ARG:HG3	2.31	0.50
1:A:48:ASN:HD21	1:A:234:THR:H	1.58	0.50
1:B:341:LEU:HD13	1:B:420:ARG:HB3	1.93	0.50
1:A:103:ARG:NH1	1:A:103:ARG:CG	2.73	0.50
1:A:34:SER:H	1:A:37:GLN:NE2	2.09	0.50
1:B:474:ILE:HG22	4:B:677:HOH:O	2.12	0.50
1:A:248:LYS:HB2	1:A:325:LEU:HA	1.95	0.49
1:B:343:LEU:N	1:B:344:PRO:CD	2.75	0.49
1:A:275:ARG:HH22	1:A:303:GLN:NE2	2.10	0.49
1:B:423:ILE:HD12	1:B:423:ILE:H	1.77	0.49
1:B:396:ARG:NH1	1:B:428:SER:OG	2.46	0.49
1:A:346:ARG:NH2	4:A:794:HOH:O	2.43	0.48
1:A:88:GLN:O	1:A:88:GLN:HG3	2.13	0.48
1:A:29:HIS:CE1	1:A:59:GLU:HA	2.49	0.48
1:B:89:ILE:CG2	1:B:423:ILE:HD13	2.44	0.48
1:B:39:GLU:OE1	1:B:324:ASP:OD2	2.33	0.47
1:B:342:ILE:HA	1:B:420:ARG:HH21	1.79	0.47
1:B:218:GLU:N	1:B:218:GLU:OE2	2.47	0.47
1:B:217:PRO:HG2	1:B:218:GLU:OE2	2.15	0.47
1:A:103:ARG:CD	4:A:1003:HOH:O	2.63	0.46
1:A:311:ILE:HG13	1:A:346:ARG:NH1	2.30	0.46
1:B:94:THR:HG21	1:B:335:PRO:HB3	1.96	0.46
1:A:209:THR:HG21	1:A:366:ILE:HD11	1.98	0.46
1:B:34:SER:H	1:B:37:GLN:NE2	2.12	0.46
1:B:413:ASP:OD1	1:B:416:ARG:HB3	2.17	0.45
1:B:459:TRP:HB3	1:B:464:HIS:CG	2.51	0.45
1:B:463:ARG:NH1	4:B:1192:HOH:O	2.49	0.45
1:A:191:LEU:HD12	4:A:1087:HOH:O	2.16	0.45
1:A:94:THR:HG21	1:A:335:PRO:HB3	1.98	0.45
1:A:39:GLU:OE1	1:A:324:ASP:OD2	2.35	0.45
1:A:89:ILE:CG2	1:A:89:ILE:O	2.64	0.44
1:A:252:LEU:H	1:A:252:LEU:HD23	1.82	0.44
1:B:250:LYS:HE3	4:B:877:HOH:O	2.16	0.44
1:B:305:ARG:NH2	4:B:1015:HOH:O	2.50	0.44
1:A:342:ILE:HG23	1:A:421:TYR:OH	2.18	0.44
1:B:81:GLU:HG2	1:B:188:ALA:CB	2.48	0.44
1:B:41:GLN:HA	1:B:41:GLN:OE1	2.18	0.44
1:A:286:VAL:HG21	1:A:304:PHE:CE1	2.48	0.43
1:B:109:GLU:OE2	4:B:817:HOH:O	2.21	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:GLN:HB2	1:A:88:GLN:HE21	1.62	0.43
1:A:116:ARG:O	1:A:120:GLU:HG3	2.18	0.43
1:B:338:VAL:HG12	1:B:424:MET:HE2	1.98	0.43
1:A:364:GLU:HG2	1:A:397:ALA:HA	2.01	0.43
1:A:27:PHE:CD2	1:A:38:PHE:HB2	2.54	0.43
1:B:136:SER:OG	4:B:870:HOH:O	2.04	0.43
1:B:305:ARG:HD3	4:B:720:HOH:O	2.18	0.43
1:B:474:ILE:HG12	1:B:474:ILE:H	1.58	0.42
1:A:393:PRO:HG2	1:A:396:ARG:CG	2.50	0.42
1:B:63:ILE:HG23	1:B:86:ILE:HG21	2.01	0.42
1:B:47:ALA:HB2	4:B:853:HOH:O	2.20	0.42
1:A:393:PRO:HG2	1:A:396:ARG:HG3	2.02	0.41
1:A:211:VAL:HB	1:A:399:ASN:HD21	1.85	0.41
1:A:29:HIS:HE1	1:A:59:GLU:HA	1.84	0.41
1:A:463:ARG:CD	4:A:790:HOH:O	2.60	0.41
1:B:41:GLN:OE1	1:B:220:PHE:HB2	2.21	0.41
1:A:332:SER:CB	1:A:339:GLU:HB2	2.51	0.41
1:B:23:ARG:HD2	1:B:131:ASP:CG	2.41	0.41
1:A:211:VAL:CB	1:A:399:ASN:HD21	2.33	0.41
1:B:345:GLU:OE1	1:B:420:ARG:NH2	2.54	0.41
1:A:182:ASN:HA	1:A:183:PRO:HD2	1.95	0.41
1:A:194:GLU:CA	4:A:738:HOH:O	2.56	0.41
1:A:34:SER:N	1:A:37:GLN:HE22	2.12	0.41
1:A:41:GLN:HA	1:A:41:GLN:OE1	2.21	0.41
1:B:104:GLU:HG2	4:B:883:HOH:O	2.20	0.41
1:B:409:ASP:OD2	1:B:411:LYS:HB2	2.21	0.40
1:A:191:LEU:O	1:B:278:ARG:HD3	2.20	0.40
1:B:350:ALA:O	1:B:353:GLU:N	2.50	0.40
1:B:86:ILE:O	1:B:89:ILE:HG12	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	462/484 (96%)	453 (98%)	9 (2%)	0	100	100
1	B	452/484 (93%)	444 (98%)	8 (2%)	0	100	100
All	All	914/968 (94%)	897 (98%)	17 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	393/407 (97%)	384 (98%)	9 (2%)	50	34
1	B	386/407 (95%)	380 (98%)	6 (2%)	62	51
All	All	779/814 (96%)	764 (98%)	15 (2%)	57	43

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	GLN
1	A	91	ASP
1	A	96	ASP
1	A	104	GLU
1	A	187	PRO
1	A	191	LEU
1	A	252	LEU
1	A	371	GLU
1	A	469	GLN
1	B	174	ILE
1	B	252	LEU
1	B	333	LYS
1	B	357	LEU
1	B	423	ILE
1	B	474	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	HIS
1	A	37	GLN
1	A	48	ASN
1	A	88	GLN
1	A	240	GLN
1	A	303	GLN
1	A	399	ASN
1	A	469	GLN
1	A	484	HIS
1	B	37	GLN
1	B	48	ASN
1	B	303	GLN
1	B	399	ASN
1	B	469	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](#)

Of 7 ligands modelled in this entry, 5 are monoatomic - leaving 2 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	$\# Z > 2$
2	1PE	A	500	-	15,15,15	1.04	1 (6%)	14,14,14	0.40	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	1PE	B	500	-	12,12,15	0.94	0	11,11,14	0.38	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	1PE	A	500	-	-	6/13/13/13	-
2	1PE	B	500	-	-	5/10/10/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	1PE	C24-C14	2.11	1.59	1.49

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	1PE	OH5-C14-C24-OH4
2	A	500	1PE	OH7-C16-C26-OH6
2	B	500	1PE	OH4-C13-C23-OH3
2	B	500	1PE	OH2-C12-C22-OH3
2	A	500	1PE	OH6-C15-C25-OH5
2	A	500	1PE	C24-C14-OH5-C25
2	A	500	1PE	OH4-C13-C23-OH3
2	B	500	1PE	C23-C13-OH4-C24
2	B	500	1PE	OH5-C14-C24-OH4
2	A	500	1PE	C16-C26-OH6-C15
2	B	500	1PE	C14-C24-OH4-C13

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	464/484 (95%)	0.03	17 (3%) 41 40	14, 23, 38, 73	0
1	B	454/484 (93%)	0.22	25 (5%) 25 24	15, 28, 51, 66	0
All	All	918/968 (94%)	0.12	42 (4%) 32 31	14, 25, 47, 73	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	189	LEU	14.2
1	A	191	LEU	9.9
1	A	188	ALA	9.0
1	A	190	GLY	6.7
1	B	193	GLU	4.8
1	B	408	ALA	4.6
1	B	407	GLY	4.4
1	A	192	PRO	3.9
1	B	419	VAL	3.5
1	A	193	GLU	3.5
1	A	484	HIS	3.4
1	B	238	VAL	3.2
1	A	104	GLU	3.1
1	B	93	GLU	3.1
1	B	349	PRO	3.1
1	A	483	GLU	2.9
1	A	423	ILE	2.9
1	B	347	LEU	2.9
1	B	438	GLY	2.9
1	B	414	LEU	2.8
1	B	409	ASP	2.8
1	B	411	LYS	2.8
1	B	353	GLU	2.7
1	A	308	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	244	TYR	2.5
1	B	192	PRO	2.5
1	B	413	ASP	2.4
1	A	278	ARG	2.3
1	A	93	GLU	2.3
1	A	28	VAL	2.3
1	A	419	VAL	2.3
1	B	374	GLY	2.2
1	B	410	GLY	2.2
1	B	172	GLU	2.2
1	B	344	PRO	2.2
1	A	415	ASP	2.2
1	B	352	LYS	2.1
1	B	241	ASP	2.1
1	B	83	GLY	2.1
1	B	412	SER	2.1
1	A	132	LEU	2.0
1	B	242	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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. 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(\AA^2)	Q<0.9
3	MG	B	2004	1/1	0.88	0.09	36,36,36,36	0
2	1PE	B	500	13/16	0.89	0.22	31,39,48,51	0
3	MG	A	2002	1/1	0.89	0.06	59,59,59,59	0
2	1PE	A	500	16/16	0.92	0.18	29,36,43,43	0
3	MG	B	2005	1/1	0.94	0.09	38,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MG	A	2001	1/1	0.98	0.05	31,31,31,31	0
3	MG	A	2003	1/1	0.98	0.09	22,22,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.