



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 06:42 am BST

PDB ID : 5DEN
Title : The First Structure of a Full-Length Mammalian Phenylalanine Hydroxylase Reveals the Architecture of an Auto-inhibited Tetramer
Authors : Arturo, E.C.; Loll, P.J.; Jaffe, E.K.
Deposited on : 2015-08-25
Resolution : 2.90 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
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

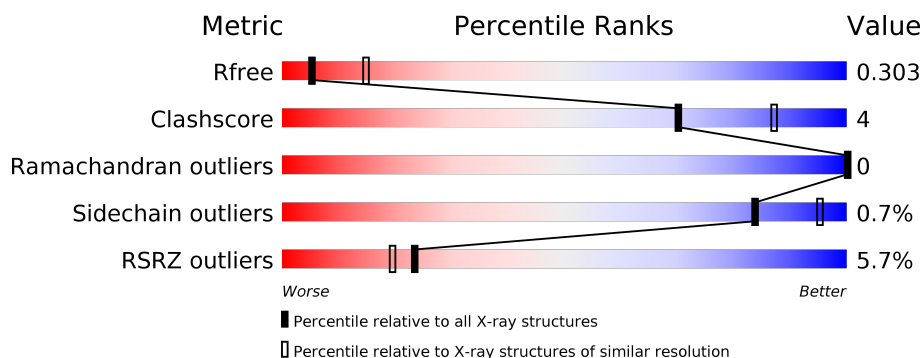
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	453	<div> <div>5%</div> <div> <div></div> <div>84%</div> <div>9%</div> <div>7%</div> </div> </div>
1	B	453	<div> <div>7%</div> <div> <div></div> <div>81%</div> <div>11%</div> <div>7%</div> </div> </div>
1	C	453	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>7%</div> </div> </div>
1	D	453	<div> <div>6%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>6%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13718 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 Phenylalanine-4-hydroxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	423	Total	C	N	O	S	0	0	0
			3435	2198	582	645	10			
1	B	420	Total	C	N	O	S	0	0	0
			3409	2181	577	641	10			
1	C	421	Total	C	N	O	S	0	0	0
			3418	2186	578	644	10			
1	D	424	Total	C	N	O	S	0	0	0
			3444	2202	583	649	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	SER	THR	conflict	UNP P04176
B	87	SER	THR	conflict	UNP P04176
C	87	SER	THR	conflict	UNP P04176
D	87	SER	THR	conflict	UNP P04176

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Fe	0	0
			1	1		
2	A	1	Total	Fe	0	0
			1	1		
2	D	1	Total	Fe	0	0
			1	1		
2	C	1	Total	Fe	0	0
			1	1		

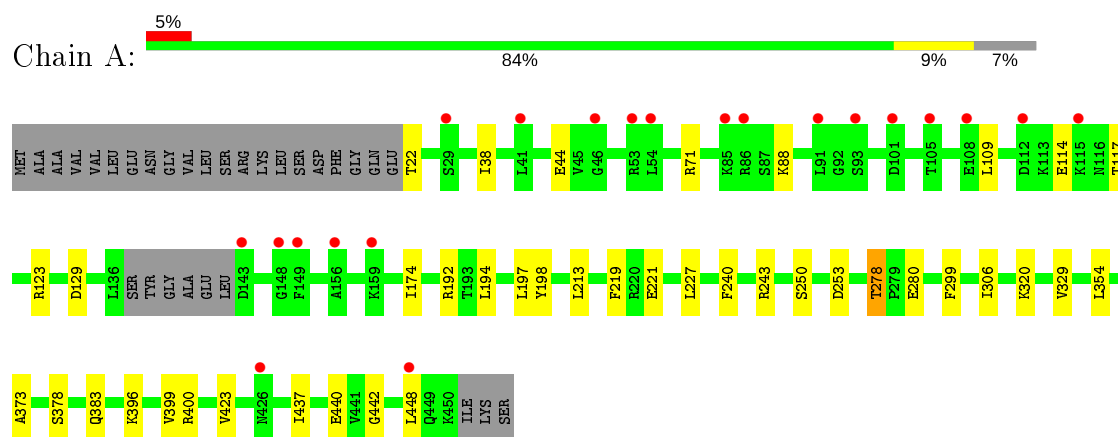
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total 2	O 2	0	0
3	B	2	Total 2	O 2	0	0
3	C	2	Total 2	O 2	0	0
3	D	2	Total 2	O 2	0	0

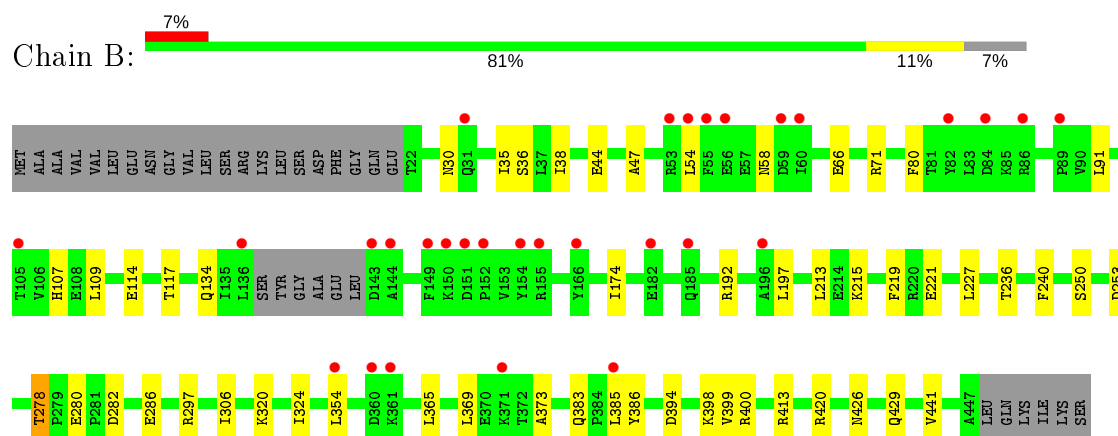
3 Residue-property plots [i](#)

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.

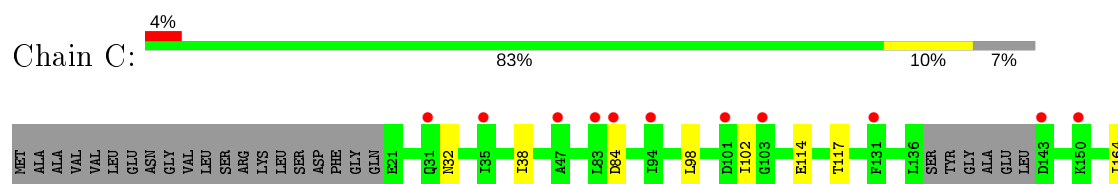
• Molecule 1: Phenylalanine-4-hydroxylase

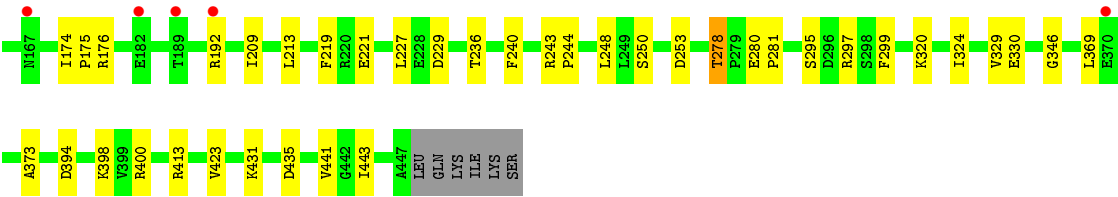


• Molecule 1: Phenylalanine-4-hydroxylase

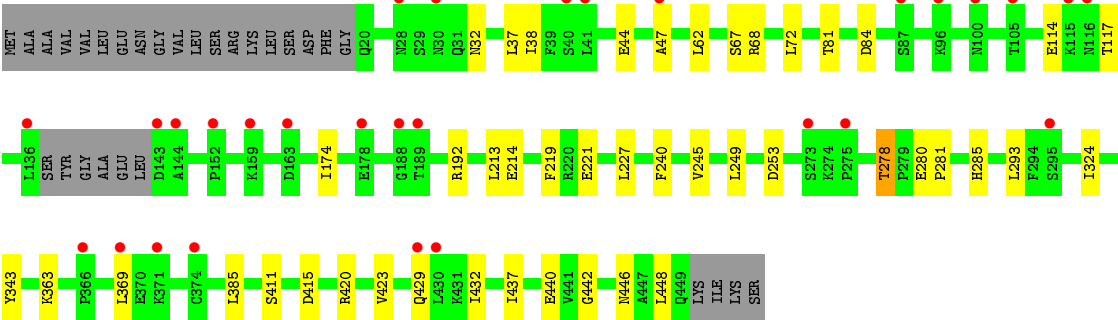
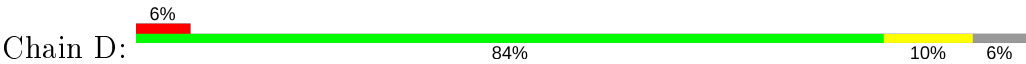


• Molecule 1: Phenylalanine-4-hydroxylase





● Molecule 1: Phenylalanine-4-hydroxylase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	113.78 Å 89.16 Å 196.81 Å 90.00° 104.53° 90.00°	Depositor
Resolution (Å)	38.10 – 2.90 38.10 – 2.90	Depositor EDS
% Data completeness (in resolution range)	90.9 (38.10-2.90) 91.0 (38.10-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 2.90 Å)	Xtriage
Refinement program	PHENIX 1.10pre_2120: ???	Depositor
R, R_{free}	0.238 , 0.303 0.238 , 0.303	Depositor DCC
R_{free} test set	1933 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	66.4	Xtriage
Anisotropy	0.565	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.017 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	13718	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.53% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.24	0/3519	0.38	0/4764
1	B	0.24	0/3493	0.38	0/4730
1	C	0.24	0/3502	0.38	0/4742
1	D	0.24	0/3528	0.38	0/4777
All	All	0.24	0/14042	0.38	0/19013

There are no bond length outliers.

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	3435	0	3366	25	1
1	B	3409	0	3334	32	0
1	C	3418	0	3340	26	0
1	D	3444	0	3367	29	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
All	All	13718	0	13407	98	1

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 4.

All (98) 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:38:ILE:HD13	1:A:423:VAL:HG21	1.77	0.67
1:D:442:GLY:O	1:D:446:ASN:ND2	2.32	0.62
1:B:38:ILE:HB	1:B:107:HIS:HB2	1.82	0.62
1:B:426:ASN:HB2	1:B:429:GLN:HG3	1.82	0.62
1:B:71:ARG:HH12	1:B:420:ARG:HB2	1.63	0.62
1:A:440:GLU:OE1	1:C:413:ARG:NH1	2.33	0.61
1:C:243:ARG:NH1	1:C:244:PRO:O	2.35	0.60
1:A:71:ARG:HG2	1:C:297:ARG:HE	1.67	0.59
1:B:413:ARG:NH1	1:D:440:GLU:OE1	2.35	0.59
1:C:394:ASP:OD2	1:C:398:LYS:NZ	2.35	0.58
1:B:278:THR:OG1	1:B:280:GLU:O	2.22	0.58
1:A:278:THR:OG1	1:A:280:GLU:O	2.21	0.57
1:B:30:ASN:ND2	1:B:134:GLN:OE1	2.30	0.57
1:C:192:ARG:NH2	1:C:221:GLU:OE1	2.36	0.56
1:D:38:ILE:HD13	1:D:423:VAL:HG21	1.86	0.56
1:B:66:GLU:HB3	1:B:80:PHE:HB2	1.87	0.56
1:D:278:THR:OG1	1:D:280:GLU:O	2.24	0.56
1:C:278:THR:OG1	1:C:280:GLU:O	2.23	0.56
1:B:250:SER:OG	1:B:253:ASP:OD2	2.23	0.56
1:A:250:SER:OG	1:A:253:ASP:OD2	2.24	0.55
1:A:299:PHE:HZ	1:A:329:VAL:HG13	1.71	0.55
1:B:192:ARG:NH2	1:B:221:GLU:OE1	2.40	0.54
1:D:293:LEU:HD11	1:D:343:TYR:HB2	1.89	0.54
1:C:38:ILE:HD13	1:C:423:VAL:HG21	1.89	0.53
1:B:36:SER:HB3	1:B:109:LEU:HD12	1.91	0.53
1:C:431:LYS:NZ	1:C:435:ASP:OD2	2.36	0.52
1:B:35:ILE:HG21	1:B:91:LEU:HD11	1.90	0.52
1:B:38:ILE:HD11	1:B:109:LEU:HD11	1.91	0.52
1:B:174:ILE:HG21	1:B:227:LEU:HB2	1.93	0.51
1:A:192:ARG:NH2	1:A:221:GLU:OE1	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:LYS:HB3	1:B:373:ALA:HB1	1.92	0.51
1:C:250:SER:OG	1:C:253:ASP:OD2	2.28	0.51
1:A:197:LEU:HD23	1:A:354:LEU:HD23	1.94	0.50
1:A:114:GLU:O	1:A:117:THR:OG1	2.27	0.50
1:D:324:ILE:HD13	1:D:369:LEU:HD13	1.92	0.50
1:C:114:GLU:O	1:C:117:THR:OG1	2.26	0.50
1:D:114:GLU:O	1:D:117:THR:OG1	2.22	0.50
1:B:394:ASP:OD1	1:B:398:LYS:NZ	2.42	0.50
1:A:88:LYS:NZ	1:A:114:GLU:OE2	2.43	0.49
1:B:297:ARG:NH2	1:D:415:ASP:OD2	2.44	0.49
1:B:400:ARG:HH11	1:D:446:ASN:HD21	1.61	0.48
1:C:324:ILE:HD13	1:C:369:LEU:HD13	1.96	0.48
1:D:281:PRO:HB3	1:D:285:HIS:HD2	1.78	0.48
1:A:320:LYS:HB3	1:A:373:ALA:HB1	1.95	0.47
1:B:413:ARG:HH12	1:D:440:GLU:CD	2.17	0.47
1:D:72:LEU:HD11	1:D:432:ILE:HG12	1.96	0.47
1:A:442:GLY:O	1:C:400:ARG:NH1	2.47	0.47
1:D:363:LYS:HE3	1:D:385:LEU:HD13	1.96	0.47
1:B:441:VAL:HG22	1:D:437:ILE:HD13	1.96	0.47
1:A:378:SER:O	1:A:383:GLN:NE2	2.42	0.47
1:B:197:LEU:HD23	1:B:354:LEU:HD23	1.96	0.47
1:C:320:LYS:HB3	1:C:373:ALA:HB1	1.96	0.47
1:B:324:ILE:HD13	1:B:369:LEU:HD13	1.96	0.47
1:B:54:LEU:O	1:B:58:ASN:ND2	2.44	0.47
1:A:129:ASP:OD1	1:A:243:ARG:NH1	2.40	0.46
1:D:245:VAL:HG21	1:D:249:LEU:HG	1.97	0.46
1:C:209:ILE:HG21	1:C:295:SER:HB2	1.97	0.46
1:A:194:LEU:HB3	1:A:198:TYR:CZ	2.51	0.46
1:D:67:SER:O	1:D:68:ARG:NH1	2.38	0.45
1:D:174:ILE:HG21	1:D:227:LEU:HB2	1.99	0.45
1:A:437:ILE:HD13	1:C:441:VAL:HG22	1.98	0.45
1:A:440:GLU:CD	1:C:413:ARG:HH12	2.20	0.45
1:A:448:LEU:HD21	1:D:448:LEU:HD21	1.99	0.45
1:B:215:LYS:HE3	1:D:62:LEU:HD13	1.98	0.44
1:C:32:ASN:O	1:C:84:ASP:HB2	2.17	0.44
1:A:213:LEU:HB3	1:A:219:PHE:CD1	2.52	0.44
1:D:32:ASN:O	1:D:84:ASP:HB2	2.18	0.44
1:D:37:LEU:H	1:D:81:THR:HG1	1.64	0.43
1:B:365:LEU:HD12	1:B:385:LEU:HD21	2.01	0.43
1:C:174:ILE:HG21	1:C:227:LEU:HB2	2.00	0.43
1:D:44:GLU:HB2	1:D:47:ALA:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:GLU:HB2	1:B:47:ALA:HB2	2.00	0.43
1:C:164:ILE:HG23	1:C:175:PRO:HG2	2.00	0.42
1:B:213:LEU:HB3	1:B:219:PHE:CD1	2.54	0.42
1:B:306:ILE:HG12	1:B:399:VAL:HG11	2.02	0.42
1:B:383:GLN:HB2	1:B:386:TYR:CZ	2.53	0.42
1:D:192:ARG:NH2	1:D:221:GLU:OE1	2.39	0.42
1:D:281:PRO:HB3	1:D:285:HIS:CD2	2.54	0.42
1:D:411:SER:OG	1:D:429:GLN:OE1	2.20	0.42
1:A:38:ILE:HD11	1:A:109:LEU:HD11	2.01	0.41
1:B:282:ASP:O	1:B:286:GLU:HG2	2.20	0.41
1:C:213:LEU:HB3	1:C:219:PHE:CD1	2.56	0.41
1:D:214:GLU:HA	1:D:219:PHE:HB2	2.02	0.41
1:D:213:LEU:HB3	1:D:219:PHE:CD1	2.54	0.41
1:A:396:LYS:HE3	1:A:400:ARG:HH12	1.85	0.41
1:C:299:PHE:HZ	1:C:329:VAL:HG13	1.86	0.41
1:C:330:GLU:O	1:C:346:GLY:N	2.50	0.41
1:A:174:ILE:HG21	1:A:227:LEU:HB2	2.02	0.41
1:A:38:ILE:HG12	1:A:123:ARG:NH1	2.35	0.41
1:B:71:ARG:NH1	1:B:420:ARG:HB2	2.31	0.41
1:C:98:LEU:HD23	1:C:102:ILE:HD12	2.02	0.41
1:D:293:LEU:HA	1:D:293:LEU:HD23	1.85	0.41
1:C:176:ARG:NH2	1:C:229:ASP:OD2	2.53	0.41
1:A:400:ARG:HD3	1:C:443:ILE:HG12	2.03	0.41
1:B:114:GLU:O	1:B:117:THR:OG1	2.31	0.40
1:B:236:THR:HG23	1:D:420:ARG:CZ	2.51	0.40
1:A:306:ILE:HG12	1:A:399:VAL:HG11	2.03	0.40
1:C:248:LEU:HD21	1:C:281:PRO:HG2	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:GLU:OE1	1:A:192:ARG:NH2[3_455]	2.18	0.02

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	419/453 (92%)	402 (96%)	17 (4%)	0	100	100
1	B	416/453 (92%)	399 (96%)	17 (4%)	0	100	100
1	C	417/453 (92%)	400 (96%)	17 (4%)	0	100	100
1	D	420/453 (93%)	401 (96%)	19 (4%)	0	100	100
All	All	1672/1812 (92%)	1602 (96%)	70 (4%)	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	375/399 (94%)	372 (99%)	3 (1%)	81	94
1	B	372/399 (93%)	370 (100%)	2 (0%)	88	96
1	C	373/399 (94%)	370 (99%)	3 (1%)	81	94
1	D	376/399 (94%)	373 (99%)	3 (1%)	81	94
All	All	1496/1596 (94%)	1485 (99%)	11 (1%)	84	95

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	THR
1	A	240	PHE
1	A	278	THR
1	B	240	PHE
1	B	278	THR
1	C	236	THR
1	C	240	PHE
1	C	278	THR

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Mol	Chain	Res	Type
1	D	240	PHE
1	D	253	ASP
1	D	278	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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	423/453 (93%)	0.39	21 (4%)	28 25	50, 75, 118, 144	0
1	B	420/453 (92%)	0.49	30 (7%)	16 12	46, 77, 112, 151	0
1	C	421/453 (92%)	0.43	16 (3%)	40 36	56, 79, 111, 142	0
1	D	424/453 (93%)	0.47	29 (6%)	17 13	49, 80, 114, 148	0
All	All	1688/1812 (93%)	0.45	96 (5%)	23 19	46, 78, 114, 151	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	150	LYS	6.8
1	D	100	ASN	4.9
1	C	189	THR	4.4
1	A	108	GLU	4.4
1	D	116	ASN	4.4
1	C	94	ILE	4.3
1	C	103	GLY	3.9
1	B	182	GLU	3.7
1	A	46	GLY	3.6
1	D	152	PRO	3.6
1	B	143	ASP	3.5
1	C	101	ASP	3.5
1	C	370	GLU	3.5
1	D	273	SER	3.4
1	A	143	ASP	3.4
1	A	105	THR	3.4
1	D	40	SER	3.4
1	D	178	GLU	3.3
1	B	136	LEU	3.2
1	C	31	GLN	3.1
1	A	156	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	144	ALA	3.0
1	A	93	SER	3.0
1	B	55	PHE	2.9
1	C	167	ASN	2.9
1	D	115	LYS	2.9
1	D	189	THR	2.9
1	B	151	ASP	2.9
1	C	84	ASP	2.8
1	A	149	PHE	2.8
1	B	59	ASP	2.7
1	C	47	ALA	2.7
1	B	152	PRO	2.7
1	B	56	GLU	2.7
1	B	31	GLN	2.7
1	A	115	LYS	2.6
1	A	101	ASP	2.6
1	B	371	LYS	2.6
1	A	91	LEU	2.6
1	B	154	TYR	2.6
1	B	105	THR	2.6
1	D	96	LYS	2.6
1	D	47	ALA	2.5
1	A	86	ARG	2.5
1	D	371	LYS	2.5
1	A	448	LEU	2.5
1	A	54	LEU	2.5
1	B	54	LEU	2.5
1	B	60	ILE	2.5
1	B	149	PHE	2.5
1	B	144	ALA	2.5
1	A	159	LYS	2.4
1	B	360	ASP	2.4
1	B	185	GLN	2.4
1	D	429	GLN	2.4
1	D	275	PRO	2.4
1	B	155	ARG	2.4
1	D	159	LYS	2.3
1	D	136	LEU	2.3
1	B	84	ASP	2.3
1	B	86	ARG	2.3
1	B	53	ARG	2.3
1	D	163	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	426	ASN	2.3
1	B	354	LEU	2.3
1	A	85	LYS	2.3
1	B	89	PRO	2.3
1	D	28	ASN	2.3
1	A	53	ARG	2.3
1	B	82	TYR	2.3
1	C	35	ILE	2.2
1	C	83	LEU	2.2
1	A	112	ASP	2.2
1	C	182	GLU	2.2
1	B	385	LEU	2.2
1	D	430	LEU	2.2
1	B	196	ALA	2.2
1	D	87	SER	2.2
1	A	148	GLY	2.1
1	D	30	ASN	2.1
1	D	366	PRO	2.1
1	D	105	THR	2.1
1	D	374	CYS	2.1
1	C	131	PHE	2.1
1	D	41	LEU	2.1
1	C	150	LYS	2.1
1	D	143	ASP	2.1
1	A	29	SER	2.1
1	A	41	LEU	2.1
1	B	166	TYR	2.1
1	D	369	LEU	2.1
1	D	295	SER	2.0
1	C	192	ARG	2.0
1	B	361	LYS	2.0
1	C	143	ASP	2.0
1	D	188	GLY	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 [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
2	FE	A	501	1/1	0.92	0.17	59,59,59,59	1
2	FE	D	501	1/1	0.93	0.18	83,83,83,83	1
2	FE	C	501	1/1	0.94	0.22	49,49,49,49	1
2	FE	B	501	1/1	0.97	0.17	44,44,44,44	1

6.5 Other polymers [i](#)

There are no such residues in this entry.