Full wwPDB X-ray Structure Validation Report (i)

Aug 21, 2023 – 08:40 PM EDT

PDB ID	:	2PAH
Title	:	TETRAMERIC HUMAN PHENYLALANINE HYDROXYLASE
Authors	:	Stevens, R.C.; Fusetti, F.; Erlandsen, H.
Deposited on	:	1998-05-26
Resolution	:	3.10 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 (1)) were used in the production of this report:

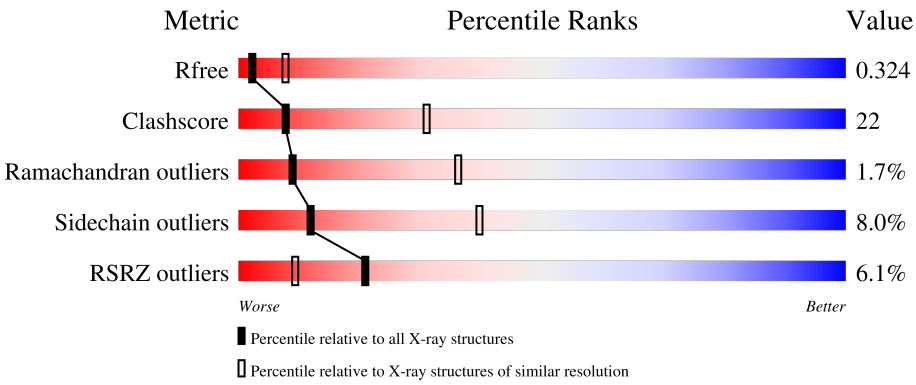
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	335	3% 51% 41% 5% •
1	B	335	9% 54% 37% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5317 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 PROTEIN (PHENYLALANINE HYDROXYLASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	329	2686	1737	448	491	10	0	0	0
1	B	322	2629	1700	439	480	10	0	0	0

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

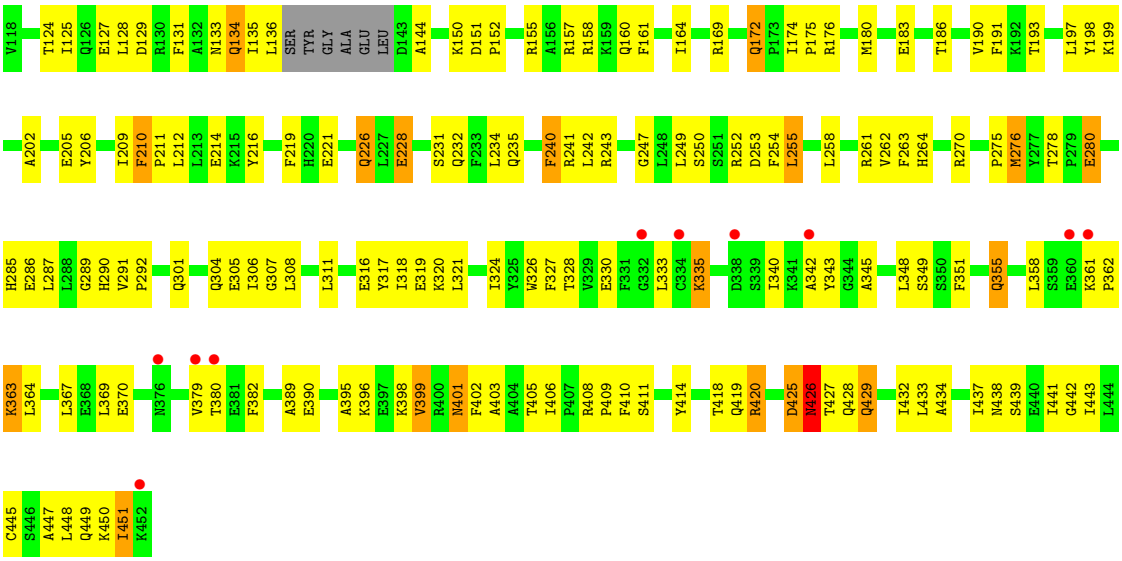
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Fe		
2	A	1	1	1	0	0
2	B	1	1	1	0	0

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.

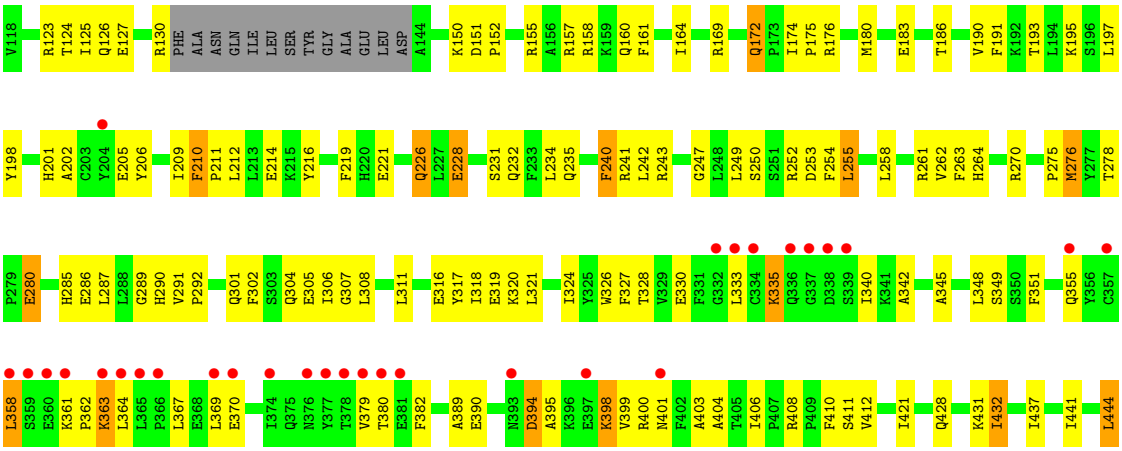
• Molecule 1: PROTEIN (PHENYLALANINE HYDROXYLASE)

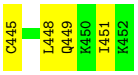
Chain A:



• Molecule 1: PROTEIN (PHENYLALANINE HYDROXYLASE)

Chain B:




C445
L448
Q449
K450
I451
K452

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 1 2	Depositor
Cell constants a, b, c, α, β, γ	119.50Å 119.50Å 126.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.10 19.92 – 3.11	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-3.10) 93.9 (19.92-3.11)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	3.70 (at 3.09Å)	Xtriage
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.251 , 0.326 0.260 , 0.324	Depositor DCC
R_{free} test set	1730 reflections (9.88%)	wwPDB-VP
Wilson B-factor (Å^2)	81.6	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 69.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	5317	wwPDB-VP
Average B, all atoms (Å^2)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% 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: 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	0.30	0/2760	0.52	0/3738
1	B	0.30	0/2702	0.52	0/3659
All	All	0.30	0/5462	0.52	0/7397

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	2686	0	2633	121	0
1	B	2629	0	2579	112	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
All	All	5317	0	5212	229	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 22.

All (229) 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:226:GLN:HE21	1:B:226:GLN:HA	1.36	0.91
1:B:369:LEU:HD13	1:B:398:LYS:HB2	1.55	0.87
1:A:226:GLN:HE21	1:A:226:GLN:HA	1.37	0.87
1:B:367:LEU:HD23	1:B:398:LYS:HE3	1.63	0.81
1:B:302:PHE:HE1	1:B:399:VAL:HG21	1.46	0.80
1:A:367:LEU:HD22	1:A:389:ALA:HB2	1.64	0.77
1:B:367:LEU:HD22	1:B:389:ALA:HB2	1.64	0.76
1:A:411:SER:HB2	1:A:429:GLN:HG3	1.70	0.74
1:B:302:PHE:CE1	1:B:399:VAL:HG21	2.21	0.74
1:B:250:SER:HB3	1:B:253:ASP:OD1	1.88	0.72
1:A:287:LEU:HD23	1:A:291:VAL:HG21	1.71	0.72
1:B:287:LEU:HD23	1:B:291:VAL:HG21	1.72	0.71
1:B:448:LEU:O	1:B:451:ILE:HG22	1.92	0.69
1:A:231:SER:HB2	1:A:242:LEU:HB2	1.74	0.69
1:A:250:SER:HB3	1:A:253:ASP:OD1	1.91	0.69
1:A:402:PHE:O	1:A:406:ILE:HG13	1.93	0.69
1:B:437:ILE:HG22	1:B:441:ILE:HD11	1.76	0.68
1:A:403:ALA:HA	1:A:406:ILE:HD12	1.75	0.68
1:B:252:ARG:NH2	1:B:318:ILE:HG13	2.08	0.67
1:B:231:SER:HB2	1:B:242:LEU:HB2	1.74	0.67
1:A:411:SER:CB	1:A:429:GLN:HG3	2.24	0.67
1:A:186:THR:HG23	1:A:275:PRO:HB3	1.75	0.67
1:A:252:ARG:NH2	1:A:318:ILE:HG13	2.09	0.67
1:A:214:GLU:HA	1:A:219:PHE:HB2	1.77	0.67
1:A:433:LEU:O	1:A:437:ILE:HG13	1.95	0.66
1:B:214:GLU:HA	1:B:219:PHE:HB2	1.76	0.66
1:B:186:THR:HG23	1:B:275:PRO:HB3	1.78	0.65
1:B:369:LEU:CD2	1:B:399:VAL:HA	2.26	0.65
1:A:335:LYS:HE2	1:A:340:ILE:HD11	1.80	0.64
1:B:335:LYS:HE2	1:B:340:ILE:HD11	1.80	0.64
1:A:129:ASP:HA	1:A:243:ARG:NH2	2.14	0.63
1:B:190:VAL:HA	1:B:193:THR:HG22	1.81	0.63
1:B:158:ARG:HH22	1:B:280:GLU:CD	2.02	0.63
1:A:124:THR:HB	1:A:127:GLU:HG3	1.80	0.63
1:A:333:LEU:HD23	1:A:342:ALA:HA	1.80	0.63
1:B:333:LEU:HD23	1:B:342:ALA:HA	1.80	0.63
1:A:425:ASP:O	1:A:428:GLN:HG2	1.97	0.63
1:A:190:VAL:HA	1:A:193:THR:HG22	1.80	0.62
1:A:158:ARG:HH22	1:A:280:GLU:CD	2.02	0.62
1:A:447:ALA:HA	1:A:450:LYS:HE2	1.80	0.62
1:B:124:THR:HB	1:B:127:GLU:HG3	1.82	0.62
1:A:361:LYS:H	1:A:362:PRO:HD2	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:451:ILE:HG23	1:B:451:ILE:O	2.01	0.61
1:B:361:LYS:H	1:B:362:PRO:HD2	1.65	0.61
1:A:151:ASP:O	1:A:155:ARG:HB2	2.00	0.61
1:A:241:ARG:HG3	1:A:262:VAL:HG22	1.82	0.61
1:B:151:ASP:O	1:B:155:ARG:HB2	2.01	0.60
1:A:428:GLN:O	1:A:432:ILE:HG12	2.01	0.60
1:A:176:ARG:NH1	1:A:226:GLN:HG3	2.16	0.60
1:A:186:THR:O	1:A:190:VAL:HG23	2.02	0.59
1:B:241:ARG:HG3	1:B:262:VAL:HG22	1.84	0.59
1:B:176:ARG:NH1	1:B:226:GLN:HG3	2.17	0.59
1:B:186:THR:O	1:B:190:VAL:HG23	2.02	0.59
1:A:193:THR:HG21	1:A:276:MET:CE	2.33	0.59
1:A:395:ALA:O	1:A:399:VAL:HG23	2.02	0.59
1:B:193:THR:HG21	1:B:276:MET:CE	2.33	0.59
1:A:286:GLU:O	1:A:291:VAL:HG23	2.02	0.58
1:A:135:ILE:O	1:A:136:LEU:HB2	2.03	0.58
1:B:193:THR:HG21	1:B:276:MET:HE2	1.85	0.58
1:B:361:LYS:H	1:B:362:PRO:CD	2.17	0.57
1:A:206:TYR:CE1	1:A:348:LEU:HD13	2.40	0.57
1:A:308:LEU:HD21	1:A:414:TYR:HB2	1.86	0.57
1:A:361:LYS:H	1:A:362:PRO:CD	2.17	0.57
1:B:286:GLU:O	1:B:291:VAL:HG23	2.03	0.57
1:B:320:LYS:O	1:B:324:ILE:HG12	2.05	0.57
1:B:206:TYR:CE1	1:B:348:LEU:HD13	2.40	0.56
1:B:228:GLU:O	1:B:232:GLN:HB2	2.06	0.56
1:A:304:GLN:HG2	1:A:308:LEU:HD11	1.87	0.56
1:A:418:THR:HB	1:A:420:ARG:HH21	1.71	0.56
1:A:190:VAL:HA	1:A:193:THR:CG2	2.35	0.56
1:B:304:GLN:HG2	1:B:308:LEU:HD11	1.87	0.55
1:B:400:ARG:O	1:B:403:ALA:HB3	2.05	0.55
1:A:320:LYS:O	1:A:324:ILE:HG12	2.06	0.55
1:B:335:LYS:HG3	1:B:390:GLU:HA	1.87	0.55
1:A:361:LYS:N	1:A:362:PRO:CD	2.70	0.55
1:B:190:VAL:HA	1:B:193:THR:CG2	2.36	0.55
1:A:228:GLU:O	1:A:232:GLN:HB2	2.08	0.54
1:A:335:LYS:HG3	1:A:390:GLU:HA	1.89	0.54
1:A:361:LYS:N	1:A:362:PRO:HD2	2.23	0.54
1:B:361:LYS:N	1:B:362:PRO:CD	2.71	0.54
1:A:169:ARG:O	1:A:172:GLN:HG2	2.07	0.54
1:A:135:ILE:HG22	1:A:136:LEU:H	1.73	0.54
1:A:408:ARG:C	1:A:410:PHE:H	2.11	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:364:LEU:H	1:A:364:LEU:HD22	1.73	0.54
1:A:438:ASN:HB2	1:B:445:CYS:SG	2.48	0.53
1:A:240:PHE:HA	1:A:261:ARG:O	2.09	0.53
1:A:193:THR:HG21	1:A:276:MET:HE2	1.90	0.53
1:A:442:GLY:O	1:A:445:CYS:HB2	2.08	0.53
1:B:326:TRP:HA	1:B:330:GLU:HB2	1.91	0.53
1:B:361:LYS:N	1:B:362:PRO:HD2	2.23	0.53
1:B:127:GLU:O	1:B:130:ARG:HG2	2.09	0.53
1:A:205:GLU:CD	1:A:205:GLU:H	2.13	0.52
1:A:212:LEU:O	1:A:216:TYR:HD1	1.92	0.52
1:A:441:ILE:HG23	1:A:442:GLY:N	2.25	0.52
1:B:306:ILE:HG23	1:B:321:LEU:HD22	1.92	0.52
1:B:212:LEU:O	1:B:216:TYR:HD1	1.92	0.52
1:A:210:PHE:HB3	1:A:211:PRO:HD3	1.91	0.52
1:B:364:LEU:HD22	1:B:364:LEU:H	1.75	0.52
1:A:198:TYR:O	1:A:202:ALA:HB3	2.10	0.51
1:A:326:TRP:HA	1:A:330:GLU:HB2	1.91	0.51
1:A:401:ASN:HD22	1:A:402:PHE:H	1.57	0.51
1:B:198:TYR:O	1:B:202:ALA:HB3	2.09	0.51
1:B:205:GLU:H	1:B:205:GLU:CD	2.12	0.51
1:B:285:HIS:HA	1:B:349:SER:CB	2.41	0.51
1:B:169:ARG:O	1:B:172:GLN:HG2	2.10	0.51
1:B:210:PHE:HB3	1:B:211:PRO:HD3	1.92	0.51
1:B:226:GLN:HA	1:B:226:GLN:NE2	2.17	0.51
1:A:252:ARG:CZ	1:A:318:ILE:HG13	2.41	0.51
1:A:135:ILE:HG22	1:A:136:LEU:N	2.26	0.50
1:B:240:PHE:HA	1:B:261:ARG:O	2.10	0.50
1:A:291:VAL:HB	1:A:292:PRO:HD3	1.93	0.50
1:A:150:LYS:O	1:A:152:PRO:HD3	2.12	0.50
1:A:425:ASP:O	1:A:427:THR:N	2.44	0.50
1:B:291:VAL:HB	1:B:292:PRO:HD3	1.94	0.50
1:B:369:LEU:HD22	1:B:398:LYS:O	2.11	0.50
1:A:306:ILE:HG23	1:A:321:LEU:HD22	1.93	0.50
1:B:191:PHE:CD2	1:B:221:GLU:HB3	2.47	0.50
1:B:252:ARG:CZ	1:B:318:ILE:HG13	2.41	0.50
1:A:191:PHE:CD2	1:A:221:GLU:HB3	2.47	0.49
1:A:451:ILE:HG23	1:A:451:ILE:O	2.12	0.49
1:A:403:ALA:HA	1:A:406:ILE:CD1	2.43	0.49
1:B:234:LEU:HD11	1:B:291:VAL:HG13	1.95	0.49
1:A:285:HIS:HA	1:A:349:SER:CB	2.42	0.49
1:A:305:GLU:HG3	1:A:396:LYS:HG3	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:448:LEU:O	1:A:451:ILE:HG22	2.13	0.49
1:A:231:SER:O	1:A:235:GLN:HG3	2.13	0.48
1:A:234:LEU:HD11	1:A:291:VAL:HG13	1.96	0.48
1:A:401:ASN:HD22	1:A:402:PHE:N	2.11	0.48
1:B:408:ARG:HH21	1:B:412:VAL:HG22	1.79	0.48
1:A:128:LEU:HD22	1:A:131:PHE:CE1	2.49	0.48
1:B:445:CYS:O	1:B:449:GLN:HG2	2.14	0.48
1:B:212:LEU:O	1:B:216:TYR:HB2	2.14	0.48
1:B:437:ILE:O	1:B:441:ILE:HG13	2.14	0.48
1:B:150:LYS:O	1:B:152:PRO:HD3	2.14	0.47
1:B:369:LEU:HD22	1:B:399:VAL:HA	1.95	0.47
1:B:123:ARG:O	1:B:421:ILE:N	2.47	0.47
1:B:197:LEU:HD11	1:B:351:PHE:CE1	2.49	0.47
1:A:304:GLN:HG2	1:A:308:LEU:CD1	2.45	0.47
1:A:425:ASP:HB3	1:A:428:GLN:CD	2.35	0.47
1:B:209:ILE:HD12	1:B:212:LEU:HD12	1.96	0.47
1:A:133:ASN:O	1:A:134:GLN:CB	2.63	0.47
1:A:209:ILE:HD12	1:A:212:LEU:HD12	1.96	0.47
1:A:414:TYR:OH	1:A:419:GLN:HA	2.15	0.47
1:B:304:GLN:HG2	1:B:308:LEU:CD1	2.45	0.47
1:A:439:SER:O	1:A:443:ILE:HD13	2.15	0.46
1:B:270:ARG:HB3	1:B:278:THR:HG23	1.96	0.46
1:B:301:GLN:O	1:B:305:GLU:HG2	2.15	0.46
1:A:197:LEU:HD11	1:A:351:PHE:CE1	2.50	0.46
1:A:270:ARG:HB3	1:A:278:THR:HG23	1.97	0.46
1:B:400:ARG:O	1:B:404:ALA:N	2.47	0.46
1:A:369:LEU:HD12	1:A:369:LEU:H	1.81	0.46
1:A:255:LEU:HA	1:A:258:LEU:HD12	1.97	0.46
1:B:395:ALA:HA	1:B:398:LYS:HE3	1.98	0.46
1:B:255:LEU:HA	1:B:258:LEU:HD12	1.98	0.46
1:A:212:LEU:O	1:A:216:TYR:HB2	2.15	0.46
1:B:124:THR:HG22	1:B:125:ILE:N	2.30	0.46
1:A:157:ARG:O	1:A:160:GLN:HB3	2.16	0.46
1:A:326:TRP:C	1:A:328:THR:H	2.20	0.46
1:B:231:SER:O	1:B:235:GLN:HG3	2.16	0.46
1:A:226:GLN:HA	1:A:226:GLN:NE2	2.18	0.45
1:A:449:GLN:HE22	1:B:431:LYS:HD3	1.81	0.45
1:B:444:LEU:HD12	1:B:444:LEU:O	2.16	0.45
1:B:399:VAL:O	1:B:403:ALA:N	2.47	0.45
1:A:124:THR:HG22	1:A:125:ILE:N	2.31	0.45
1:A:301:GLN:O	1:A:305:GLU:HG2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:ARG:O	1:A:410:PHE:N	2.48	0.45
1:B:164:ILE:CG2	1:B:175:PRO:HG2	2.46	0.45
1:B:382:PHE:N	1:B:382:PHE:CD1	2.84	0.45
1:B:379:VAL:HG13	1:B:380:THR:HG23	1.99	0.45
1:A:164:ILE:CG2	1:A:175:PRO:HG2	2.46	0.45
1:A:190:VAL:CA	1:A:193:THR:HG22	2.46	0.45
1:B:390:GLU:HB2	1:B:394:ASP:CG	2.38	0.45
1:A:379:VAL:HG13	1:A:380:THR:HG23	1.99	0.45
1:B:307:GLY:O	1:B:311:LEU:HG	2.17	0.45
1:B:157:ARG:O	1:B:160:GLN:HB3	2.16	0.45
1:B:243:ARG:HD3	1:B:264:HIS:CE1	2.52	0.44
1:B:226:GLN:HE21	1:B:226:GLN:CA	2.16	0.44
1:A:402:PHE:O	1:A:402:PHE:CD2	2.71	0.44
1:B:326:TRP:C	1:B:328:THR:H	2.19	0.44
1:A:328:THR:O	1:A:343:TYR:CE2	2.70	0.44
1:A:161:PHE:O	1:A:164:ILE:HB	2.18	0.44
1:A:243:ARG:HD3	1:A:264:HIS:CE1	2.53	0.44
1:B:316:GLU:HG2	1:B:317:TYR:N	2.33	0.44
1:B:242:LEU:HD23	1:B:263:PHE:HB3	2.00	0.43
1:A:242:LEU:HD23	1:A:263:PHE:HB3	2.00	0.43
1:A:369:LEU:HD12	1:A:369:LEU:N	2.34	0.43
1:A:382:PHE:N	1:A:382:PHE:CD1	2.84	0.43
1:A:441:ILE:HG21	1:B:441:ILE:HD12	2.00	0.43
1:B:369:LEU:H	1:B:369:LEU:HD12	1.81	0.43
1:A:316:GLU:HG2	1:A:317:TYR:N	2.33	0.43
1:B:190:VAL:CA	1:B:193:THR:HG22	2.47	0.43
1:A:307:GLY:O	1:A:311:LEU:HG	2.18	0.43
1:A:428:GLN:HE21	1:A:428:GLN:HB3	1.72	0.43
1:A:363:LYS:HB3	1:A:363:LYS:HE2	1.83	0.43
1:B:124:THR:HG21	1:B:126:GLN:OE1	2.19	0.43
1:B:174:ILE:N	1:B:228:GLU:HG2	2.34	0.43
1:A:316:GLU:HG2	1:A:317:TYR:H	1.84	0.42
1:B:369:LEU:HD21	1:B:399:VAL:HG22	2.01	0.42
1:A:434:ALA:HB1	1:B:445:CYS:SG	2.59	0.42
1:B:316:GLU:HG2	1:B:317:TYR:H	1.84	0.42
1:B:201:HIS:ND1	1:B:358:LEU:HG	2.35	0.42
1:B:369:LEU:HD12	1:B:369:LEU:N	2.34	0.42
1:A:174:ILE:N	1:A:228:GLU:HG2	2.34	0.42
1:B:161:PHE:O	1:B:164:ILE:HB	2.19	0.42
1:B:395:ALA:O	1:B:399:VAL:HG23	2.20	0.42
1:A:402:PHE:HD2	1:A:405:THR:OG1	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:379:VAL:HG13	1:B:380:THR:N	2.35	0.41
1:A:425:ASP:O	1:A:426:ASN:C	2.58	0.41
1:B:130:ARG:HD2	1:B:130:ARG:N	2.34	0.41
1:B:403:ALA:HA	1:B:406:ILE:HG12	2.02	0.41
1:B:180:MET:O	1:B:183:GLU:HB2	2.20	0.41
1:B:206:TYR:CD1	1:B:348:LEU:HD13	2.56	0.41
1:B:198:TYR:HD1	1:B:202:ALA:HB2	1.85	0.41
1:A:379:VAL:HG13	1:A:380:THR:N	2.35	0.41
1:A:290:HIS:CD2	1:A:345:ALA:HB1	2.56	0.41
1:B:191:PHE:O	1:B:195:LYS:HB2	2.21	0.41
1:B:212:LEU:O	1:B:216:TYR:CD1	2.74	0.41
1:A:206:TYR:CD1	1:A:348:LEU:HD13	2.56	0.40
1:A:355:GLN:H	1:A:355:GLN:HG2	1.65	0.40
1:B:432:ILE:H	1:B:432:ILE:HG12	1.42	0.40
1:B:316:GLU:O	1:B:320:LYS:HG3	2.21	0.40
1:A:395:ALA:HA	1:A:398:LYS:HG3	2.03	0.40
1:B:290:HIS:CD2	1:B:345:ALA:HB1	2.57	0.40
1:A:180:MET:O	1:A:183:GLU:HB2	2.22	0.40
1:A:335:LYS:HB2	1:A:390:GLU:O	2.22	0.40
1:A:382:PHE:N	1:A:382:PHE:HD1	2.19	0.40
1:B:363:LYS:HE2	1:B:363:LYS:HB3	1.83	0.40
1:A:199:LYS:HE3	1:A:199:LYS:HB3	1.98	0.40
1:A:398:LYS:H	1:A:398:LYS:HG2	1.68	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	325/335 (97%)	291 (90%)	26 (8%)	8 (2%)	5 27
1	B	318/335 (95%)	288 (91%)	27 (8%)	3 (1%)	17 52

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	643/670 (96%)	579 (90%)	53 (8%)	11 (2%)	9 36

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	426	ASN
1	A	134	GLN
1	A	144	ALA
1	A	289	GLY
1	B	289	GLY
1	A	247	GLY
1	A	327	PHE
1	B	247	GLY
1	B	327	PHE
1	A	409	PRO
1	A	451	ILE

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	290/294 (99%)	268 (92%)	22 (8%)	13 41
1	B	284/294 (97%)	260 (92%)	24 (8%)	10 37
All	All	574/588 (98%)	528 (92%)	46 (8%)	12 40

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	172	GLN
1	A	210	PHE
1	A	226	GLN
1	A	228	GLU
1	A	240	PHE
1	A	249	LEU
1	A	254	PHE

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Mol	Chain	Res	Type
1	A	255	LEU
1	A	276	MET
1	A	280	GLU
1	A	319	GLU
1	A	335	LYS
1	A	355	GLN
1	A	358	LEU
1	A	363	LYS
1	A	370	GLU
1	A	399	VAL
1	A	401	ASN
1	A	420	ARG
1	A	425	ASP
1	A	426	ASN
1	A	429	GLN
1	B	172	GLN
1	B	210	PHE
1	B	226	GLN
1	B	228	GLU
1	B	240	PHE
1	B	249	LEU
1	B	254	PHE
1	B	255	LEU
1	B	276	MET
1	B	280	GLU
1	B	319	GLU
1	B	335	LYS
1	B	355	GLN
1	B	358	LEU
1	B	363	LYS
1	B	370	GLU
1	B	394	ASP
1	B	398	LYS
1	B	401	ASN
1	B	410	PHE
1	B	411	SER
1	B	428	GLN
1	B	432	ILE
1	B	444	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	167	ASN
1	A	170	HIS
1	A	264	HIS
1	A	383	GLN
1	A	401	ASN
1	A	419	GLN
1	A	449	GLN
1	B	167	ASN
1	B	170	HIS
1	B	264	HIS
1	B	383	GLN
1	B	401	ASN
1	B	426	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)

Of 2 ligands modelled in this entry, 2 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 (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	329/335 (98%)	-0.05	10 (3%) 50 27	17, 66, 98, 100	0
1	B	322/335 (96%)	0.34	30 (9%) 8 3	18, 79, 100, 100	0
All	All	651/670 (97%)	0.15	40 (6%) 21 9	17, 72, 99, 100	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	338	ASP	6.9
1	B	337	GLY	6.2
1	B	393	ASN	5.3
1	B	355	GLN	5.0
1	B	332	GLY	4.3
1	B	334	CYS	3.8
1	B	374	ILE	3.7
1	A	379	VAL	3.5
1	A	332	GLY	3.4
1	B	379	VAL	3.4
1	A	338	ASP	3.3
1	B	360	GLU	3.3
1	A	380	THR	3.0
1	A	452	LYS	2.9
1	B	357	CYS	2.9
1	A	342	ALA	2.9
1	A	361	LYS	2.9
1	B	333	LEU	2.8
1	B	364	LEU	2.7
1	B	397	GLU	2.7
1	B	401	ASN	2.7
1	B	336	GLN	2.7
1	A	360	GLU	2.6
1	B	359	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	334	CYS	2.6
1	B	361	LYS	2.6
1	B	381	GLU	2.4
1	A	376	ASN	2.4
1	B	358	LEU	2.4
1	B	363	LYS	2.3
1	B	365	LEU	2.3
1	B	377	TYR	2.2
1	B	380	THR	2.2
1	B	369	LEU	2.1
1	B	339	SER	2.1
1	B	378	THR	2.1
1	B	370	GLU	2.0
1	B	366	PRO	2.0
1	B	376	ASN	2.0
1	B	204	TYR	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
2	FE	B	454	1/1	0.55	0.12	91,91,91,91	0
2	FE	A	453	1/1	0.73	0.07	84,84,84,84	0

6.5 Other polymers (i)

There are no such residues in this entry.