



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:31 PM UTC

PDB ID : 3RNG / pdb_00003rng
Title : Structure of the Toluene/o-Xylene Monooxygenase Hydroxylase
T201S/W167E Double Mutant
Authors : Gucinski, G.; Song, W.J.; Lippard, S.J.; Sazinsky, M.H.
Deposited on : 2011-04-22
Resolution : 2.81 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

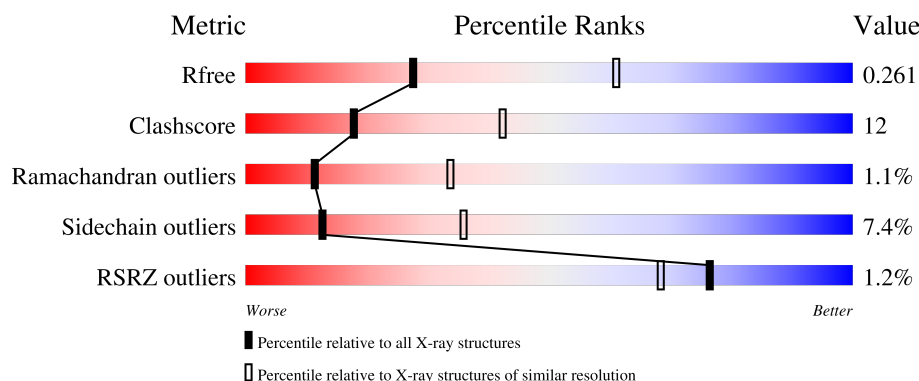
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

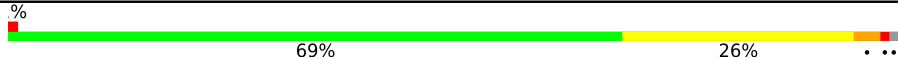
The reported resolution of this entry is 2.81 Å.

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}	180053	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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 $\leq 5\%$. 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	498	 2% 69% 26% ...
2	B	330	 2% 74% 19% . .
3	C	86	 2% 58% 28% 9% . .

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7373 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 Toluene o-xylene monooxygenase component.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	491	Total	C	N	O	S	0	0	0
			4006	2555	672	753	26			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	167	GLU	TRP	engineered mutation	UNP Q6IV66
A	201	SER	THR	engineered mutation	UNP Q6IV66
A	445	LYS	GLU	engineered mutation	UNP Q6IV66

- Molecule 2 is a protein called Toluene o-xylene monooxygenase component.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	322	Total	C	N	O	S	0	0	0
			2641	1674	467	490	10			

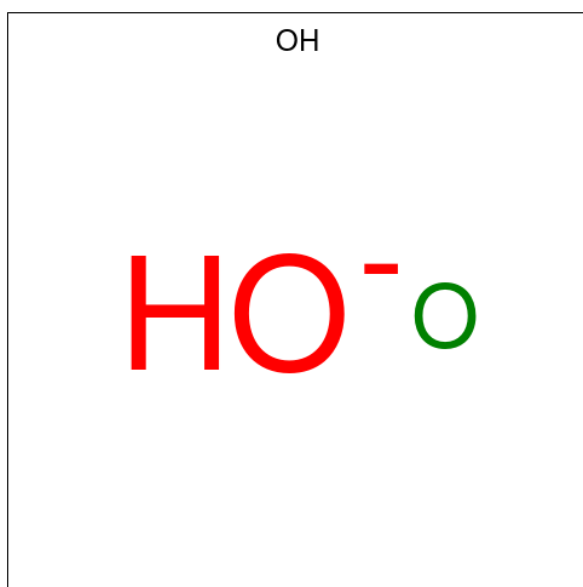
- Molecule 3 is a protein called Toluene o-xylene monooxygenase component.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	83	Total	C	N	O	S	0	0	0
			676	425	120	126	5			

- Molecule 4 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Fe	0	0
			2	2		

- Molecule 5 is HYDROXIDE ION (CCD ID: OH) (formula: HO).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	O	0	0
			1	1		
5	A	1	Total	O	0	0
			1	1		

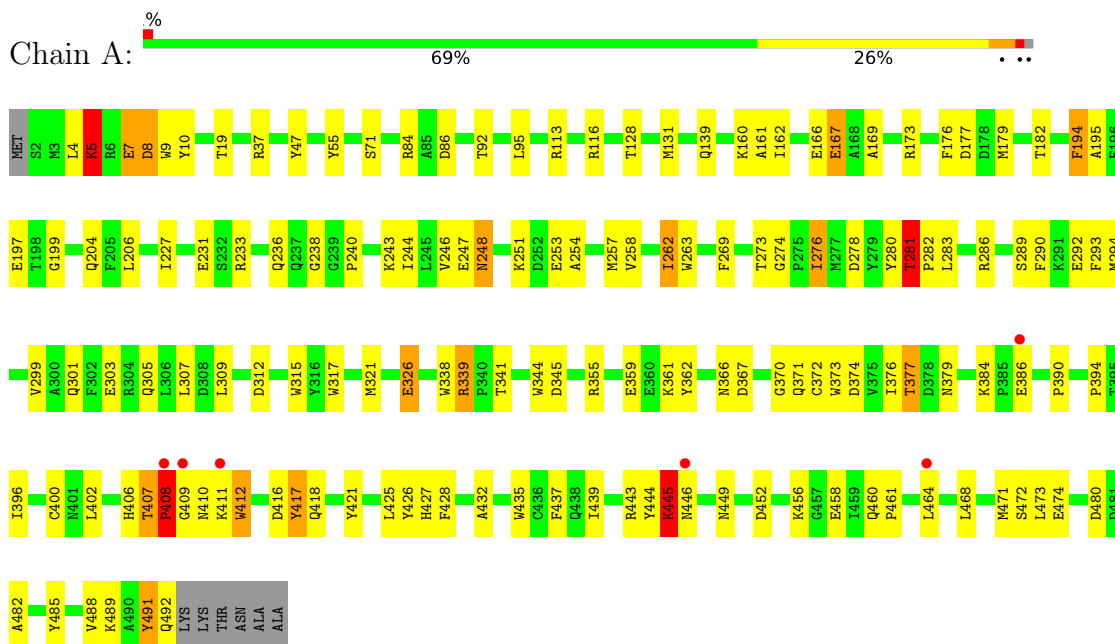
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	26	Total	O	0	0
			26	26		
6	B	18	Total	O	0	0
			18	18		
6	C	2	Total	O	0	0
			2	2		

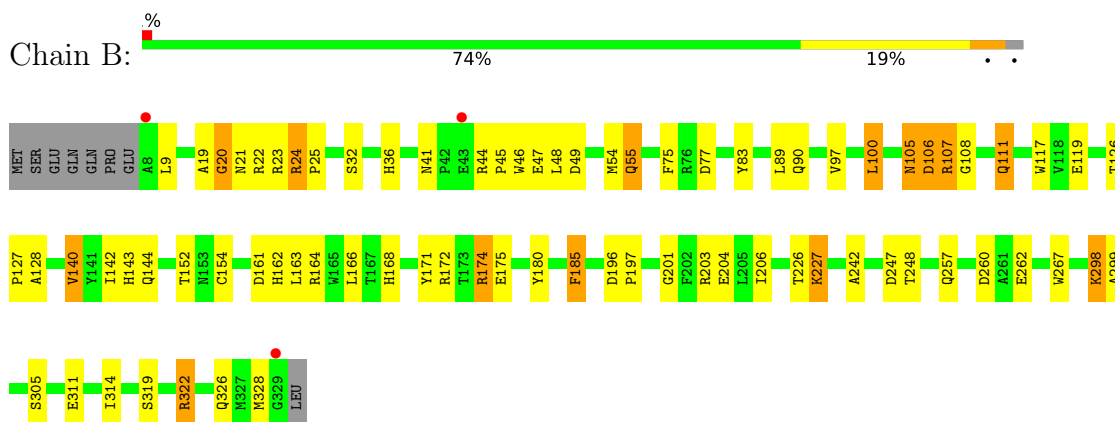
3 Residue-property plots

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: Toluene o-xylene monooxygenase component



- Molecule 2: Toluene o-xylene monooxygenase component



- Molecule 3: Toluene o-xylene monooxygenase component





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	182.63Å 182.63Å 68.28Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.66 – 2.81 45.66 – 2.81	Depositor EDS
% Data completeness (in resolution range)	96.0 (45.66-2.81) 96.2 (45.66-2.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.02 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.192 , 0.261 0.196 , 0.261	Depositor DCC
R_{free} test set	1568 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	55.9	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 31.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7373	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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: OH, 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	0.91	0/4128	1.13	16/5609 (0.3%)
2	B	1.00	0/2713	1.14	7/3688 (0.2%)
3	C	0.83	0/690	1.25	6/934 (0.6%)
All	All	0.94	0/7531	1.15	29/10231 (0.3%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	326	GLU	N-CA-C	12.90	129.69	113.43
1	A	408	PRO	CB-CA-C	-8.00	98.37	111.56
2	B	227	LYS	CA-C-N	-7.79	110.71	119.28
2	B	227	LYS	C-N-CA	-7.79	110.71	119.28
3	C	60	THR	N-CA-C	7.65	121.96	110.64
3	C	4	PHE	CB-CA-C	7.53	120.36	109.45
1	A	326	GLU	CB-CA-C	-7.07	97.77	109.65
1	A	281	THR	N-CA-C	-6.89	101.07	109.83
1	A	84	ARG	CB-CA-C	6.80	120.43	109.27
2	B	185	PHE	CB-CA-C	-6.76	100.31	109.71
1	A	386	GLU	N-CA-C	-6.55	105.28	113.28
1	A	281	THR	CB-CA-C	6.51	119.65	109.42
2	B	55	GLN	N-CA-C	-6.24	104.40	111.14
3	C	12	ARG	CB-CA-C	-6.09	101.30	111.23
3	C	77	GLU	N-CA-C	6.07	118.85	110.35
1	A	86	ASP	CA-C-N	-5.99	113.64	119.87
1	A	86	ASP	C-N-CA	-5.99	113.64	119.87
2	B	75	PHE	N-CA-C	5.96	117.98	110.24
3	C	38	SER	N-CA-C	5.94	121.08	113.18
1	A	339	ARG	N-CA-C	5.67	121.33	113.45
1	A	379	ASN	N-CA-C	5.54	118.04	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	41	ARG	N-CA-C	-5.50	105.62	112.93
2	B	180	TYR	CA-C-N	5.33	125.00	119.56
2	B	180	TYR	C-N-CA	5.33	125.00	119.56
1	A	417	TYR	N-CA-C	-5.33	97.23	107.57
1	A	8	ASP	N-CA-C	5.23	117.88	111.82
1	A	227	ILE	CB-CA-C	-5.14	105.30	112.04
1	A	194	PHE	N-CA-C	5.09	116.68	111.03
1	A	5	LYS	N-CA-C	-5.09	102.23	110.32

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	4006	0	3756	112	0
2	B	2641	0	2537	70	0
3	C	676	0	667	25	0
4	A	2	0	0	0	0
5	A	2	0	0	0	0
6	A	26	0	0	1	0
6	B	18	0	0	0	0
6	C	2	0	0	0	0
All	All	7373	0	6960	175	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (175) 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:446:ASN:HB2	2:B:49:ASP:OD1	1.36	1.21
1:A:446:ASN:CB	2:B:49:ASP:OD1	2.15	0.94
2:B:162:HIS:HE1	2:B:227:LYS:HZ2	1.02	0.92
1:A:139:GLN:HE22	2:B:83:TYR:H	1.15	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:GLU:OE1	1:A:366:ASN:ND2	2.05	0.89
2:B:162:HIS:CE1	2:B:227:LYS:HZ2	1.90	0.87
1:A:418:GLN:HE22	3:C:78:THR:H	1.28	0.80
2:B:162:HIS:HE1	2:B:227:LYS:NZ	1.80	0.79
1:A:427:HIS:HE1	3:C:76:THR:HG23	1.45	0.79
2:B:47:GLU:O	2:B:48:LEU:HG	1.83	0.79
1:A:446:ASN:HB2	2:B:49:ASP:CG	2.07	0.79
3:C:75:PRO:O	3:C:76:THR:HG22	1.84	0.77
3:C:54:ARG:HG3	3:C:80:ASP:HB2	1.66	0.75
1:A:411:LYS:O	1:A:412:TRP:C	2.30	0.75
2:B:126:THR:OG1	2:B:127:PRO:HD3	1.86	0.75
1:A:92:THR:HG23	1:A:276:ILE:HG12	1.67	0.75
1:A:367:ASP:HB3	1:A:410:ASN:HD22	1.55	0.71
1:A:305:GLN:HE21	1:A:305:GLN:HA	1.56	0.70
1:A:197:GLU:HG2	1:A:231:GLU:OE2	1.91	0.70
2:B:168:HIS:HD2	2:B:257:GLN:HE21	1.41	0.69
1:A:7:GLU:H	1:A:7:GLU:CD	2.01	0.68
1:A:47:TYR:CE1	1:A:240:PRO:HB2	2.29	0.67
1:A:253:GLU:OE1	1:A:253:GLU:N	2.24	0.67
1:A:408:PRO:HA	1:A:412:TRP:HD1	1.59	0.67
1:A:244:ILE:O	1:A:248:ASN:HB2	1.95	0.67
1:A:485:TYR:O	1:A:488:VAL:HB	1.95	0.65
1:A:37:ARG:NH2	1:A:257:MET:HE1	2.11	0.65
1:A:113:ARG:HH11	2:B:144:GLN:HE21	1.45	0.64
1:A:139:GLN:HE22	2:B:83:TYR:N	1.90	0.64
1:A:282:PRO:O	1:A:286:ARG:HG3	1.97	0.64
3:C:28:ASP:OD1	3:C:64:ARG:HB3	1.99	0.63
2:B:77:ASP:HB2	2:B:154:CYS:SG	2.38	0.63
2:B:105:ASN:O	2:B:107:ARG:N	2.32	0.63
1:A:95:LEU:HD13	1:A:344:TRP:CH2	2.35	0.62
1:A:446:ASN:CB	2:B:49:ASP:CG	2.69	0.62
1:A:139:GLN:NE2	2:B:83:TYR:H	1.94	0.62
1:A:278:ASP:OD2	1:A:362:TYR:OH	2.13	0.61
1:A:408:PRO:HB3	3:C:13:ASP:C	2.25	0.61
1:A:372:CYS:O	1:A:376:ILE:HG12	2.01	0.61
1:A:8:ASP:O	2:B:174:ARG:HD2	2.02	0.60
1:A:326:GLU:HG3	1:A:407:THR:OG1	2.02	0.60
1:A:317:TRP:O	1:A:321:MET:HG2	2.02	0.59
2:B:305:SER:HB3	2:B:314:ILE:HD11	1.83	0.59
1:A:9:TRP:HB3	2:B:174:ARG:HG3	1.83	0.59
1:A:367:ASP:HB3	1:A:410:ASN:ND2	2.17	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:GLU:OE1	1:A:407:THR:HG23	2.02	0.58
1:A:194:PHE:HB2	1:A:238:GLY:HA3	1.85	0.58
1:A:345:ASP:O	1:A:482:ALA:HB2	2.03	0.57
2:B:163:LEU:HD12	2:B:166:LEU:HD23	1.85	0.57
1:A:446:ASN:ND2	2:B:44:ARG:HH22	2.03	0.57
3:C:30:VAL:HG12	3:C:31:ALA:N	2.20	0.56
1:A:416:ASP:OD2	1:A:427:HIS:HD2	1.89	0.56
1:A:446:ASN:HD22	2:B:44:ARG:HH22	1.54	0.56
1:A:243:LYS:O	1:A:247:GLU:HG2	2.06	0.56
1:A:373:TRP:O	1:A:377:THR:OG1	2.24	0.55
2:B:168:HIS:CD2	2:B:257:GLN:HE21	2.22	0.55
2:B:111:GLN:HE21	2:B:111:GLN:H	1.54	0.55
2:B:161:ASP:O	2:B:164:ARG:HB3	2.05	0.55
2:B:196:ASP:O	2:B:197:PRO:C	2.49	0.55
1:A:246:VAL:C	1:A:248:ASN:H	2.14	0.55
1:A:408:PRO:HA	1:A:412:TRP:CD1	2.41	0.55
2:B:201:GLY:HA3	2:B:299:ALA:HA	1.90	0.54
2:B:204:GLU:OE1	2:B:298:LYS:NZ	2.34	0.54
3:C:60:THR:O	3:C:60:THR:HG22	2.06	0.54
1:A:366:ASN:HA	1:A:370:GLY:HA3	1.89	0.53
1:A:446:ASN:HB3	2:B:49:ASP:HA	1.90	0.53
2:B:117:TRP:NE1	2:B:247:ASP:HB2	2.24	0.53
1:A:258:VAL:O	1:A:262:ILE:HG13	2.08	0.53
1:A:263:TRP:CE2	1:A:432:ALA:HB3	2.44	0.53
3:C:63:PRO:HG2	3:C:66:MET:HE3	1.89	0.53
1:A:425:LEU:HD23	3:C:76:THR:HG22	1.90	0.52
1:A:19:THR:O	2:B:203:ARG:NH2	2.43	0.52
2:B:105:ASN:HD22	2:B:106:ASP:N	2.07	0.52
1:A:491:TYR:CD2	1:A:491:TYR:N	2.77	0.52
3:C:77:GLU:HG3	3:C:78:THR:N	2.26	0.51
1:A:471:MET:HE3	1:A:473:LEU:HD12	1.93	0.51
2:B:154:CYS:HB3	2:B:267:TRP:CE2	2.46	0.51
2:B:107:ARG:O	2:B:107:ARG:HG3	2.11	0.51
1:A:4:LEU:HD11	2:B:175:GLU:HG2	1.93	0.50
2:B:105:ASN:C	2:B:107:ARG:H	2.19	0.50
3:C:35:ALA:HB1	3:C:39:ILE:HG13	1.94	0.50
1:A:169:ALA:HB1	1:A:173:ARG:NH1	2.27	0.50
1:A:464:LEU:O	1:A:468:LEU:HG	2.12	0.50
2:B:19:ALA:O	2:B:20:GLY:O	2.30	0.49
3:C:23:THR:O	3:C:68:VAL:HB	2.12	0.49
1:A:233:ARG:O	1:A:236:GLN:NE2	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:ARG:HD2	1:A:480:ASP:HA	1.94	0.49
1:A:491:TYR:O	1:A:492:GLN:HB2	2.12	0.49
2:B:105:ASN:O	2:B:106:ASP:C	2.54	0.49
2:B:117:TRP:CZ2	2:B:242:ALA:HA	2.47	0.49
1:A:411:LYS:O	1:A:412:TRP:O	2.30	0.49
2:B:126:THR:C	2:B:128:ALA:H	2.20	0.49
1:A:418:GLN:NE2	3:C:78:THR:H	2.04	0.48
1:A:417:TYR:O	1:A:428:PHE:HB2	2.12	0.48
1:A:246:VAL:C	1:A:248:ASN:N	2.70	0.48
1:A:338:TRP:CE2	1:A:394:PRO:HD3	2.49	0.48
2:B:105:ASN:C	2:B:107:ARG:N	2.67	0.48
1:A:444:TYR:O	1:A:445:LYS:C	2.55	0.48
1:A:443:ARG:HG2	1:A:444:TYR:CZ	2.49	0.47
1:A:400:CYS:SG	1:A:402:LEU:HB2	2.54	0.47
2:B:126:THR:C	2:B:128:ALA:N	2.72	0.47
1:A:294:MET:O	1:A:299:VAL:HG23	2.15	0.47
2:B:105:ASN:O	2:B:108:GLY:N	2.28	0.47
1:A:9:TRP:CE2	2:B:171:TYR:HD1	2.33	0.47
1:A:301:GLN:O	1:A:305:GLN:HG2	2.14	0.46
2:B:164:ARG:NH2	2:B:260:ASP:OD2	2.46	0.46
1:A:315:TRP:HZ2	1:A:435:TRP:CE3	2.33	0.46
2:B:36:HIS:HE1	2:B:152:THR:OG1	1.98	0.46
2:B:154:CYS:HB3	2:B:267:TRP:CD2	2.51	0.46
1:A:449:ASN:ND2	1:A:452:ASP:OD2	2.49	0.46
2:B:47:GLU:C	2:B:48:LEU:HG	2.40	0.46
1:A:305:GLN:HA	1:A:305:GLN:NE2	2.26	0.46
1:A:427:HIS:CE1	3:C:76:THR:HG23	2.37	0.46
1:A:116:ARG:HB2	2:B:140:VAL:HG13	1.98	0.46
1:A:162:ILE:O	1:A:173:ARG:NH1	2.49	0.46
1:A:408:PRO:HB3	3:C:13:ASP:O	2.16	0.46
1:A:243:LYS:O	1:A:247:GLU:CG	2.65	0.45
1:A:421:TYR:HB3	1:A:426:TYR:HE2	1.79	0.45
2:B:143:HIS:CD2	2:B:143:HIS:C	2.94	0.45
1:A:9:TRP:O	1:A:10:TYR:C	2.59	0.45
3:C:21:VAL:HG11	3:C:30:VAL:HG22	1.99	0.45
2:B:142:ILE:O	2:B:143:HIS:C	2.58	0.45
1:A:195:ALA:O	1:A:199:GLY:HA3	2.17	0.44
1:A:421:TYR:HB3	1:A:426:TYR:CE2	2.52	0.44
1:A:5:LYS:HG3	1:A:8:ASP:OD2	2.16	0.44
1:A:254:ALA:O	1:A:258:VAL:HG23	2.17	0.44
1:A:372:CYS:HB3	3:C:42:ARG:HG2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:437:PHE:C	1:A:437:PHE:CD2	2.95	0.44
2:B:206:ILE:HD13	2:B:226:THR:HG21	1.99	0.44
2:B:46:TRP:O	2:B:48:LEU:N	2.49	0.44
3:C:4:PHE:HA	3:C:5:PRO:HD2	1.92	0.44
1:A:128:THR:O	1:A:131:MET:HB3	2.17	0.44
1:A:456:LYS:HE2	1:A:456:LYS:HB3	1.67	0.44
1:A:176:PHE:O	1:A:177:ASP:C	2.59	0.44
1:A:338:TRP:CZ2	1:A:394:PRO:HD3	2.52	0.44
1:A:113:ARG:HH11	2:B:144:GLN:NE2	2.13	0.43
3:C:7:MET:HG2	3:C:76:THR:O	2.17	0.43
1:A:182:THR:HA	2:B:54:MET:HG2	2.00	0.43
3:C:75:PRO:O	3:C:76:THR:CG2	2.61	0.43
1:A:446:ASN:HD22	2:B:44:ARG:NH2	2.16	0.43
3:C:74:ARG:O	3:C:77:GLU:HB2	2.18	0.43
1:A:160:LYS:O	1:A:161:ALA:C	2.61	0.43
1:A:371:GLN:O	1:A:374:ASP:HB2	2.19	0.43
2:B:90:GLN:HA	2:B:90:GLN:OE1	2.18	0.43
2:B:322:ARG:HB3	2:B:326:GLN:HE21	1.84	0.43
1:A:408:PRO:HA	1:A:412:TRP:HA	2.00	0.42
1:A:204:GLN:HB2	1:A:269:PHE:HZ	1.83	0.42
2:B:19:ALA:O	2:B:20:GLY:C	2.61	0.42
1:A:55:TYR:CE1	2:B:164:ARG:HD2	2.55	0.42
2:B:172:ARG:HA	2:B:172:ARG:HD2	1.80	0.42
3:C:46:GLN:C	3:C:48:GLU:H	2.28	0.42
2:B:126:THR:N	2:B:127:PRO:CD	2.83	0.41
2:B:45:PRO:HD2	2:B:55:GLN:OE1	2.20	0.41
2:B:105:ASN:C	2:B:105:ASN:ND2	2.79	0.41
1:A:289:SER:OG	1:A:292:GLU:HG3	2.20	0.41
3:C:74:ARG:O	3:C:75:PRO:C	2.63	0.41
1:A:281:THR:HG23	6:A:519:HOH:O	2.19	0.41
1:A:359:GLU:HB2	1:A:366:ASN:HD21	1.86	0.41
1:A:274:GLY:HA2	1:A:290:PHE:CD1	2.56	0.41
2:B:97:VAL:O	2:B:100:LEU:HB2	2.20	0.41
1:A:167:GLU:H	1:A:167:GLU:HG3	1.27	0.41
1:A:289:SER:HB3	1:A:361:LYS:HD2	2.03	0.41
1:A:293:PHE:CD2	1:A:293:PHE:C	2.98	0.41
1:A:446:ASN:ND2	2:B:44:ARG:NH2	2.69	0.41
1:A:446:ASN:N	2:B:49:ASP:OD1	2.50	0.41
1:A:460:GLN:HA	1:A:461:PRO:C	2.45	0.41
1:A:416:ASP:OD1	1:A:416:ASP:C	2.64	0.40
1:A:166:GLU:OE2	1:A:341:THR:O	2.38	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:355:ARG:HE	1:A:355:ARG:HB3	1.74	0.40
1:A:179:MET:HA	1:A:179:MET:HE2	2.03	0.40
1:A:341:THR:HA	1:A:473:LEU:HD13	2.04	0.40
1:A:409:GLY:HA2	3:C:12:ARG:HB3	2.03	0.40
2:B:126:THR:HG21	2:B:185:PHE:CD1	2.56	0.40
2:B:322:ARG:O	2:B:326:GLN:HG3	2.21	0.40
2:B:24:ARG:HA	2:B:25:PRO:HD3	1.91	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	489/498 (98%)	444 (91%)	40 (8%)	5 (1%)	12	36
2	B	320/330 (97%)	303 (95%)	14 (4%)	3 (1%)	14	38
3	C	81/86 (94%)	68 (84%)	11 (14%)	2 (2%)	4	15
All	All	890/914 (97%)	815 (92%)	65 (7%)	10 (1%)	11	34

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	408	PRO
1	A	412	TRP
1	A	445	LYS
2	B	20	GLY
2	B	106	ASP
1	A	489	LYS
2	B	22	ARG
3	C	35	ALA
1	A	390	PRO
3	C	68	VAL

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	415/422 (98%)	387 (93%)	28 (7%)	15	40
2	B	274/282 (97%)	253 (92%)	21 (8%)	12	34
3	C	77/79 (98%)	69 (90%)	8 (10%)	7	21
All	All	766/783 (98%)	709 (93%)	57 (7%)	13	36

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	7	GLU
1	A	71	SER
1	A	167	GLU
1	A	206	LEU
1	A	248	ASN
1	A	251	LYS
1	A	262	ILE
1	A	273	THR
1	A	276	ILE
1	A	280	TYR
1	A	281	THR
1	A	283	LEU
1	A	303	GLU
1	A	307	LEU
1	A	309	LEU
1	A	312	ASP
1	A	377	THR
1	A	384	LYS
1	A	396	ILE
1	A	406	HIS
1	A	407	THR
1	A	439	ILE
1	A	445	LYS
1	A	458	GLU
1	A	472	SER

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Mol	Chain	Res	Type
1	A	474	GLU
1	A	491	TYR
2	B	9	LEU
2	B	21	ASN
2	B	23	ARG
2	B	24	ARG
2	B	32	SER
2	B	41	ASN
2	B	89	LEU
2	B	100	LEU
2	B	105	ASN
2	B	107	ARG
2	B	111	GLN
2	B	119	GLU
2	B	140	VAL
2	B	174	ARG
2	B	248	THR
2	B	262	GLU
2	B	298	LYS
2	B	311	GLU
2	B	319	SER
2	B	322	ARG
2	B	328	MET
3	C	3	THR
3	C	23	THR
3	C	26	THR
3	C	30	VAL
3	C	50	ILE
3	C	57	GLU
3	C	60	THR
3	C	76	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	GLN
1	A	141	GLN
1	A	237	GLN
1	A	248	ASN
1	A	256	GLN
1	A	288	GLN
1	A	305	GLN

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Mol	Chain	Res	Type
1	A	379	ASN
1	A	401	ASN
1	A	418	GLN
1	A	427	HIS
1	A	446	ASN
1	A	447	HIS
2	B	21	ASN
2	B	34	ASN
2	B	36	HIS
2	B	98	GLN
2	B	105	ASN
2	B	111	GLN
2	B	144	GLN
2	B	153	ASN
2	B	162	HIS
2	B	168	HIS
2	B	223	ASN
2	B	237	GLN
2	B	326	GLN
3	C	40	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic and 2 are modelled with single atom - 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 [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	491/498 (98%)	-0.25	6 (1%) 76 68	32, 51, 73, 83	0
2	B	322/330 (97%)	-0.45	3 (0%) 81 74	27, 43, 63, 88	0
3	C	83/86 (96%)	0.58	2 (2%) 59 49	57, 75, 89, 95	0
All	All	896/914 (98%)	-0.25	11 (1%) 76 68	27, 49, 78, 95	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	446	ASN	3.3
2	B	43	GLU	3.2
2	B	329	GLY	3.2
1	A	409	GLY	3.1
1	A	464	LEU	2.5
1	A	408	PRO	2.4
1	A	386	GLU	2.4
3	C	66	MET	2.4
2	B	8	ALA	2.3
1	A	411	LYS	2.1
3	C	84	MET	2.1

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 oligosaccharides 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
5	OH	A	502	1/1	0.96	0.11	39,39,39,39	0
5	OH	A	501	1/1	0.99	0.04	36,36,36,36	0
4	FE	A	499	1/1	1.00	0.02	42,42,42,42	0
4	FE	A	500	1/1	1.00	0.01	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.