



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 02:37 PM GMT

PDB ID : 3RNB
Title : Structure of the Toluene/o-Xylene Monooxygenase Hydroxylase
T201S/F176W Double Mutant
Authors : Gucinski, G.; Song, W.J.; Lippard, S.J.; Sazinsky, M.H.
Deposited on : 2011-04-22
Resolution : 2.64 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

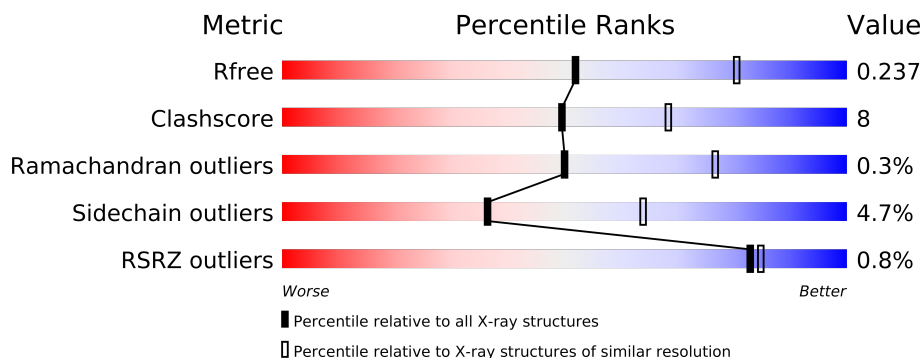
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.64 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}	66092	2393 (2.68-2.60)
Clashscore	79885	2915 (2.68-2.60)
Ramachandran outliers	78287	2865 (2.68-2.60)
Sidechain outliers	78261	2865 (2.68-2.60)
RSRZ outliers	66119	2393 (2.68-2.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	498	
2	B	330	
3	C	86	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
6	SO4	A	502	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7526 atoms, of which 0 are hydrogen and 0 are deuterium.

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	1	0
			4024	2568	676	754	26			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	176	TRP	PHE	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	323	Total	C	N	O	S	0	0	0
			2649	1680	468	491	10			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	85	Total	C	N	O	S	0	0	0
			686	431	122	128	5			

- Molecule 4 is HYDROXIDE ION (three-letter code: OH) (formula: HO).

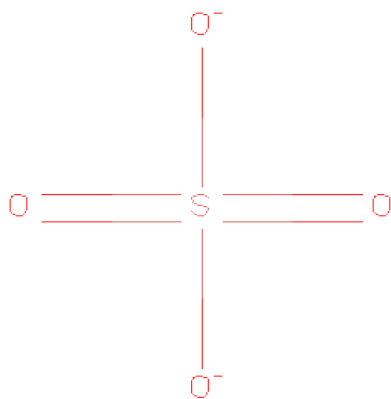


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O 1 1	0	0

- Molecule 5 is FE (III) ION (three-letter code: FE) (formula: Fe).

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

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is water.

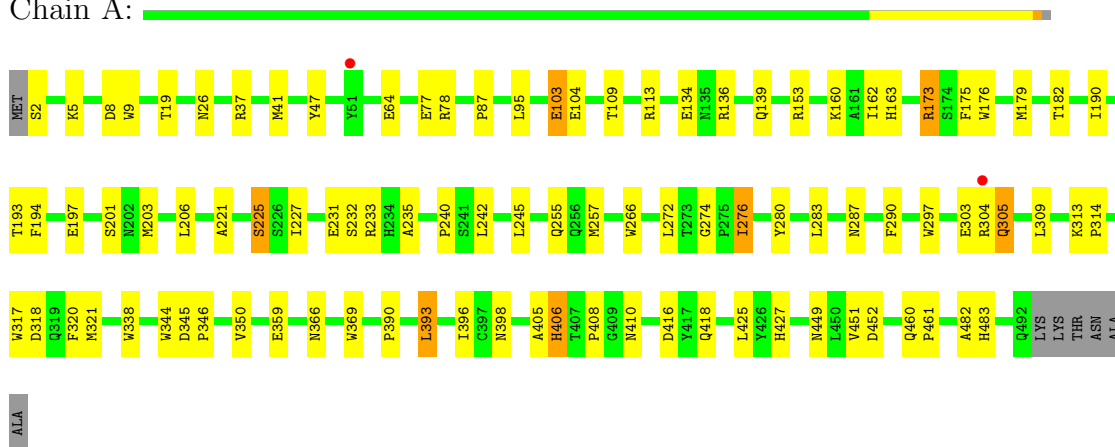
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	78	Total	O	0	0
			78	78		
7	B	65	Total	O	0	0
			65	65		
7	C	6	Total	O	0	0
			6	6		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

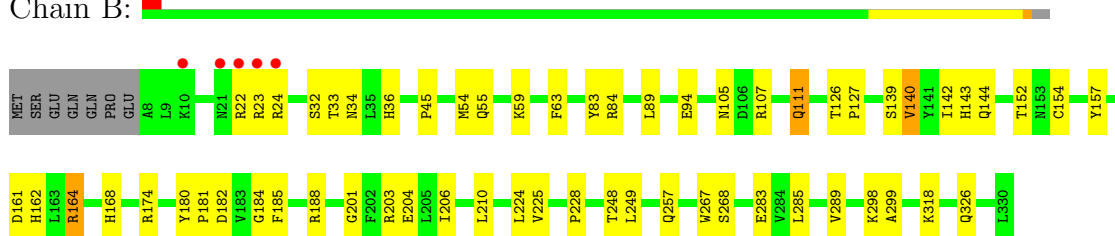
- Molecule 1: Toluene o-xylene monooxygenase component

Chain A:



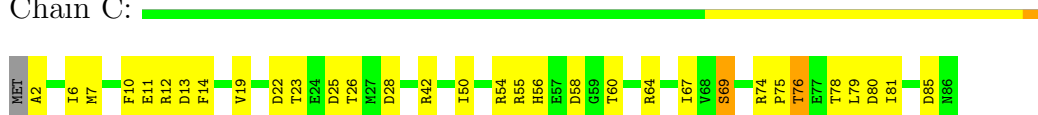
- Molecule 2: Toluene o-xylene monooxygenase component

Chain B:



- Molecule 3: Toluene o-xylene monooxygenase component

Chain C:



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	183.16Å 183.16Å 68.08Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.99 – 2.64 44.99 – 2.64	Depositor EDS
% Data completeness (in resolution range)	99.4 (44.99-2.64) 99.4 (44.99-2.64)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.17 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.175 , 0.235 0.183 , 0.237	Depositor DCC
R_{free} test set	1916 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 21.8	EDS
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 38378 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7526	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.80% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: OH, SO4, 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	0.92	0/4149	0.87	3/5640 (0.1%)
2	B	0.97	4/2721 (0.1%)	0.88	2/3699 (0.1%)
3	C	0.80	0/700	0.90	0/948
All	All	0.93	4/7570 (0.1%)	0.88	5/10287 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	283	GLU	CG-CD	5.87	1.60	1.51
2	B	182	ASP	CB-CG	5.81	1.64	1.51
2	B	157	TYR	CE2-CZ	5.54	1.45	1.38
2	B	140	VAL	CB-CG1	5.30	1.64	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	393	LEU	CA-CB-CG	7.17	131.78	115.30
1	A	103	GLU	OE1-CD-OE2	-5.94	116.17	123.30
1	A	173	ARG	NE-CZ-NH2	-5.46	117.57	120.30
2	B	84	ARG	NE-CZ-NH1	5.18	122.89	120.30
2	B	164	ARG	NE-CZ-NH1	-5.09	117.75	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens

added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4024	0	3771	61	0
2	B	2649	0	2548	42	0
3	C	686	0	674	26	0
4	A	1	0	0	1	0
5	A	2	0	0	0	0
6	A	10	0	0	1	0
6	B	5	0	0	1	0
7	A	78	0	0	1	0
7	B	65	0	0	2	0
7	C	6	0	0	0	0
All	All	7526	0	6993	109	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (109) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:369:TRP:CH2	1:A:406:HIS:O	2.12	1.01
3:C:75:PRO:O	3:C:76:THR:HB	1.61	0.98
1:A:427:HIS:HE1	3:C:76:THR:HG23	1.29	0.97
1:A:427:HIS:HE1	3:C:76:THR:CG2	1.82	0.92
1:A:8:ASP:O	2:B:174:ARG:HD2	1.71	0.90
1:A:427:HIS:CE1	3:C:76:THR:HG23	2.08	0.89
2:B:111:GLN:H	2:B:111:GLN:HE21	1.23	0.87
1:A:425:LEU:HD23	3:C:76:THR:HG22	1.59	0.85
1:A:9:TRP:HB3	2:B:174:ARG:HG3	1.61	0.81
2:B:168:HIS:HD2	2:B:257:GLN:HE21	1.27	0.80
1:A:139:GLN:HE22	2:B:83:TYR:H	1.27	0.77
2:B:126:THR:OG1	2:B:127:PRO:HD3	1.85	0.77
1:A:416:ASP:OD2	1:A:427:HIS:HD2	1.72	0.73
2:B:126:THR:HG21	2:B:185:PHE:HD1	1.55	0.72
1:A:203:MET:HG2	1:A:297:TRP:HB3	1.74	0.70
3:C:75:PRO:O	3:C:76:THR:CB	2.33	0.70
1:A:26:ASN:HB2	6:A:502:SO4:O1	1.93	0.68
1:A:136:ARG:HD2	2:B:83:TYR:CE1	2.30	0.66
1:A:255:GLN:HB3	7:A:536:HOH:O	1.96	0.65
2:B:318:LYS:NZ	6:B:331:SO4:O4	2.29	0.64
1:A:9:TRP:CB	2:B:174:ARG:HG3	2.28	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:47:TYR:CE1	1:A:240:PRO:HB2	2.34	0.63
1:A:418:GLN:HE22	3:C:78:THR:H	1.47	0.62
1:A:95:LEU:HD23	1:A:276:ILE:HD11	1.82	0.62
3:C:54:ARG:HG3	3:C:80:ASP:HB2	1.82	0.60
2:B:168:HIS:CD2	2:B:257:GLN:HE21	2.13	0.60
2:B:126:THR:HG21	2:B:185:PHE:CD1	2.36	0.60
1:A:398:ASN:HD22	1:A:427:HIS:H	1.49	0.60
1:A:2:SER:N	2:B:105:ASN:HD22	1.99	0.59
2:B:201:GLY:HA3	2:B:299:ALA:HA	1.83	0.59
3:C:11:GLU:HG2	3:C:12:ARG:HG3	1.86	0.58
1:A:318:ASP:O	1:A:321:MET:HB2	2.05	0.57
3:C:28:ASP:OD1	3:C:64:ARG:HB3	2.05	0.57
1:A:314:PRO:HD2	1:A:317:TRP:CE3	2.41	0.56
1:A:95:LEU:HD11	1:A:272:LEU:HD22	1.88	0.56
1:A:427:HIS:CE1	3:C:76:THR:CG2	2.74	0.55
2:B:224:LEU:O	2:B:228:PRO:HG2	2.06	0.55
2:B:204:GLU:OE1	2:B:298:LYS:NZ	2.39	0.55
1:A:221:ALA:O	1:A:225:SER:HB3	2.07	0.55
1:A:344:TRP:O	1:A:346:PRO:HD3	2.07	0.54
1:A:369:TRP:HH2	1:A:406:HIS:O	1.83	0.54
1:A:139:GLN:NE2	2:B:83:TYR:H	2.03	0.54
1:A:109:THR:HG23	2:B:144:GLN:HB2	1.88	0.54
1:A:19:THR:O	2:B:203:ARG:NH2	2.41	0.54
1:A:193:THR:HA	1:A:197:GLU:OE1	2.08	0.53
1:A:338:TRP:CD1	1:A:390:PRO:HG3	2.43	0.53
1:A:113:ARG:HH11	2:B:144:GLN:HE21	1.57	0.53
2:B:33:THR:HB	2:B:34:ASN:HD22	1.73	0.52
3:C:6:ILE:HD13	3:C:79:LEU:HD12	1.92	0.52
1:A:398:ASN:ND2	1:A:427:HIS:H	2.08	0.52
1:A:460:GLN:HA	1:A:461:PRO:C	2.30	0.51
1:A:369:TRP:CZ2	1:A:406:HIS:O	2.61	0.51
2:B:36:HIS:HE1	2:B:152:THR:OG1	1.93	0.51
2:B:111:GLN:N	2:B:111:GLN:HE21	2.00	0.51
1:A:410:ASN:OD1	3:C:12:ARG:NH2	2.44	0.50
1:A:416:ASP:H	3:C:56:HIS:CE1	2.28	0.50
2:B:326:GLN:HB3	7:B:370:HOH:O	2.11	0.49
1:A:232:SER:HA	1:A:235:ALA:HB3	1.93	0.49
1:A:182:THR:HA	2:B:54:MET:HG2	1.95	0.49
1:A:303:GLU:HG2	1:A:313:LYS:HE2	1.94	0.49
1:A:345:ASP:HB3	1:A:482:ALA:HA	1.95	0.49
1:A:87:PRO:HG2	1:A:483:HIS:CD2	2.48	0.48
1:A:197:GLU:HG2	1:A:231:GLU:OE2	2.15	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:175:PHE:CE1	1:A:179:MET:HG3	2.51	0.46
1:A:37:ARG:NH2	1:A:257:MET:HE1	2.31	0.46
1:A:103:GLU:HG3	1:A:176:TRP:CD1	2.50	0.46
2:B:184:GLY:HA3	2:B:188:ARG:HD3	1.97	0.46
2:B:180:TYR:N	2:B:181:PRO:HD3	2.31	0.46
2:B:94:GLU:OE1	2:B:164:ARG:NE	2.45	0.45
2:B:143:HIS:CD2	2:B:143:HIS:C	2.90	0.45
3:C:2:ALA:O	3:C:23:THR:N	2.48	0.45
1:A:396:ILE:HG12	1:A:451:VAL:HG23	1.97	0.45
3:C:55:ARG:HD2	3:C:58:ASP:OD1	2.17	0.45
2:B:249:LEU:C	2:B:249:LEU:HD23	2.37	0.45
1:A:359:GLU:HB2	1:A:366:ASN:ND2	2.32	0.45
2:B:45:PRO:HD2	2:B:55:GLN:OE1	2.16	0.45
2:B:210:LEU:HD23	2:B:210:LEU:HA	1.74	0.44
1:A:37:ARG:NH2	1:A:257:MET:CE	2.80	0.44
2:B:161:ASP:O	2:B:164:ARG:HB3	2.17	0.44
1:A:190:ILE:O	1:A:194:PHE:HB3	2.18	0.44
3:C:74:ARG:O	3:C:75:PRO:C	2.56	0.44
2:B:111:GLN:H	2:B:111:GLN:NE2	2.03	0.44
1:A:449:ASN:O	1:A:452:ASP:HB2	2.17	0.43
1:A:104:GLU:O	1:A:134:GLU:HB3	2.18	0.43
2:B:59:LYS:HA	2:B:63:PHE:HD2	1.83	0.43
3:C:10:PHE:CE1	3:C:81:ILE:HG21	2.54	0.43
1:A:162:ILE:O	1:A:173:ARG:NH1	2.52	0.43
3:C:54:ARG:HB2	3:C:60:THR:O	2.19	0.42
1:A:190:ILE:HD12	1:A:242:LEU:HG	2.00	0.42
2:B:285:LEU:O	2:B:289:VAL:HG23	2.20	0.42
1:A:305:GLN:HA	1:A:305:GLN:HE21	1.84	0.42
2:B:142:ILE:O	2:B:143:HIS:C	2.58	0.42
3:C:7:MET:HG2	3:C:76:THR:O	2.20	0.41
2:B:298:LYS:HB3	2:B:298:LYS:HE2	1.85	0.41
1:A:160:LYS:O	1:A:163:HIS:HB2	2.20	0.41
2:B:154:CYS:HB3	2:B:267:TRP:CE2	2.56	0.41
3:C:25:ASP:O	3:C:67:ILE:HA	2.21	0.41
3:C:56:HIS:HD2	3:C:80:ASP:OD1	2.04	0.41
2:B:225:VAL:C	2:B:228:PRO:HD2	2.41	0.41
1:A:274:GLY:HA2	1:A:290:PHE:CD1	2.56	0.41
1:A:231:GLU:OE1	4:A:499:OH:O	2.38	0.41
3:C:2:ALA:O	3:C:22:ASP:HA	2.20	0.41
2:B:139:SER:OG	2:B:162:HIS:HD2	2.04	0.41
1:A:405:ALA:HA	3:C:14:PHE:O	2.20	0.41
3:C:13:ASP:OD1	3:C:42:ARG:HD2	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:188:ARG:NH1	7:B:360:HOH:O	2.53	0.40
1:A:266:TRP:CD2	1:A:320:PHE:HE1	2.38	0.40
1:A:245:LEU:HD23	1:A:245:LEU:HA	1.88	0.40
3:C:6:ILE:HG13	3:C:19:VAL:O	2.20	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	490/498 (98%)	460 (94%)	28 (6%)	2 (0%)	43	70
2	B	321/330 (97%)	315 (98%)	6 (2%)	0	100	100
3	C	83/86 (96%)	75 (90%)	7 (8%)	1 (1%)	19	37
All	All	894/914 (98%)	850 (95%)	41 (5%)	3 (0%)	50	77

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	406	HIS
3	C	69	SER
1	A	408	PRO

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	417/422 (99%)	397 (95%)	20 (5%)	35	62
2	B	275/282 (98%)	264 (96%)	11 (4%)	42	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	77/79 (98%)	72 (94%)	5 (6%)	24	45
All	All	769/783 (98%)	733 (95%)	36 (5%)	36	64

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	41	MET
1	A	64	GLU
1	A	77	GLU
1	A	78	ARG
1	A	153	ARG
1	A	201	SER
1	A	206	LEU
1	A	225	SER
1	A	227	ILE
1	A	233	ARG
1	A	276	ILE
1	A	280	TYR
1	A	283	LEU
1	A	287	ASN
1	A	304	ARG
1	A	305	GLN
1	A	309	LEU
1	A	350	VAL
1	A	393	LEU
2	B	22	ARG
2	B	23	ARG
2	B	24	ARG
2	B	32	SER
2	B	89	LEU
2	B	107	ARG
2	B	111	GLN
2	B	140	VAL
2	B	206	ILE
2	B	248	THR
2	B	268	SER
3	C	26	THR
3	C	50	ILE
3	C	69	SER
3	C	76	THR
3	C	85	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	GLN
1	A	237	GLN
1	A	248	ASN
1	A	305	GLN
1	A	379	ASN
1	A	398	ASN
1	A	418	GLN
1	A	427	HIS
1	A	483	HIS
2	B	34	ASN
2	B	36	HIS
2	B	87	ASN
2	B	111	GLN
2	B	144	GLN
2	B	153	ASN
2	B	162	HIS
2	B	168	HIS
2	B	326	GLN
3	C	56	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 6 ligands modelled in this entry, 1 is modelled with single atom and 2 are monoatomic - leaving 3 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	$\# Z > 2$
6	SO4	A	502	-	4,4,4	0.39	0	6,6,6	0.71	0
6	SO4	A	503	-	4,4,4	0.35	0	6,6,6	0.38	0
6	SO4	B	331	-	4,4,4	0.45	0	6,6,6	0.25	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
6	SO4	A	502	-	-	0/0/0/0	0/0/0/0
6	SO4	A	503	-	-	0/0/0/0	0/0/0/0
6	SO4	B	331	-	-	0/0/0/0	0/0/0/0

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.

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	491/498 (98%)	-0.23	2 (0%) 90 91	29, 45, 60, 77	0
2	B	323/330 (97%)	-0.35	5 (1%) 70 70	31, 41, 60, 86	0
3	C	85/86 (98%)	-0.39	0 100 100	40, 53, 65, 74	0
All	All	899/914 (98%)	-0.29	7 (0%) 83 85	29, 45, 61, 86	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	23	ARG	3.4
2	B	22	ARG	3.2
2	B	10	LYS	2.4
2	B	21	ASN	2.3
1	A	304	ARG	2.3
1	A	51	TYR	2.1
2	B	24	ARG	2.1

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 ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	SO4	A	502	5/5	0.20	2.24	42,42,46,46	5
6	SO4	A	503	5/5	0.16	-0.29	49,51,52,52	5
4	OH	A	499	1/1	0.18	-0.64	49,49,49,49	0
5	FE	A	501	1/1	0.17	-0.94	43,43,43,43	0
5	FE	A	500	1/1	0.16	-1.25	38,38,38,38	0
6	SO4	B	331	5/5	0.11	-6.27	47,48,49,49	5

6.5 Other polymers ⓘ

There are no such residues in this entry.