Full wwPDB X-ray Structure Validation Report

Feb 1, 2016 – 12:40 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 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

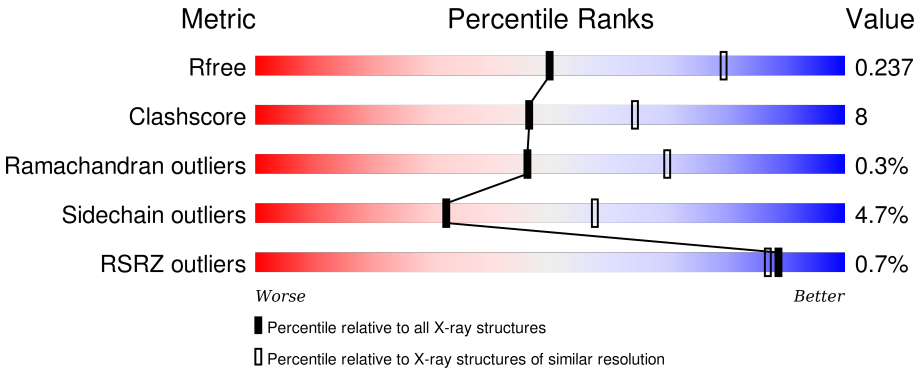
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 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}	91344	3377 (2.68-2.60)
Clashscore	102246	3781 (2.68-2.60)
Ramachandran outliers	100387	3722 (2.68-2.60)
Sidechain outliers	100360	3722 (2.68-2.60)
RSRZ outliers	91569	3388 (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. 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	498	80% 18%
2	B	330	% 80% 17%
3	C	86	62% 35%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7526 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	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

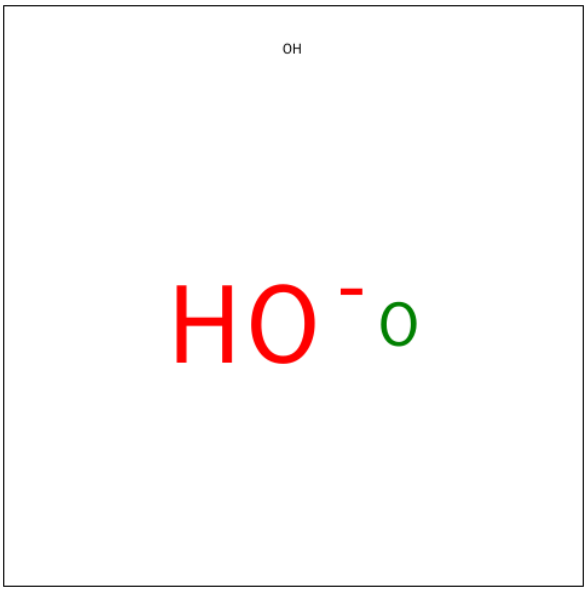
- Molecule 2 is a protein called Toluene o-xylene monooxygenase component.

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).

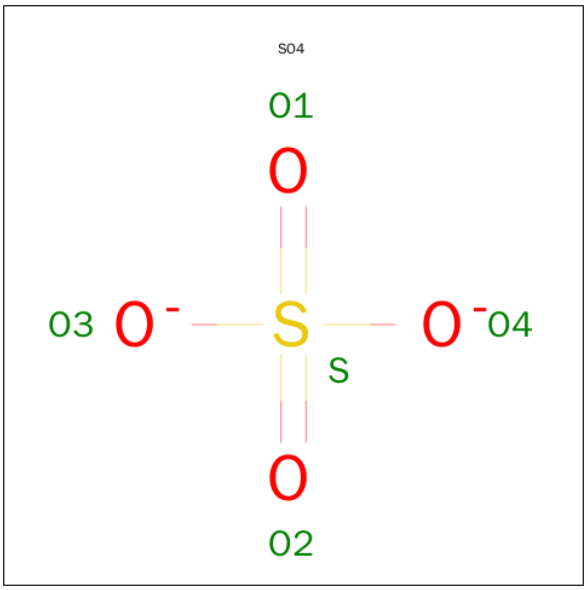


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O 1 1	0	0

- 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: O_4S).



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

- Molecule 7 is water.

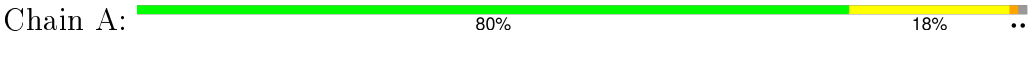
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	78	Total O 78 78	0	0
7	B	65	Total O 65 65	0	0
7	C	6	Total O 6 6	0	0

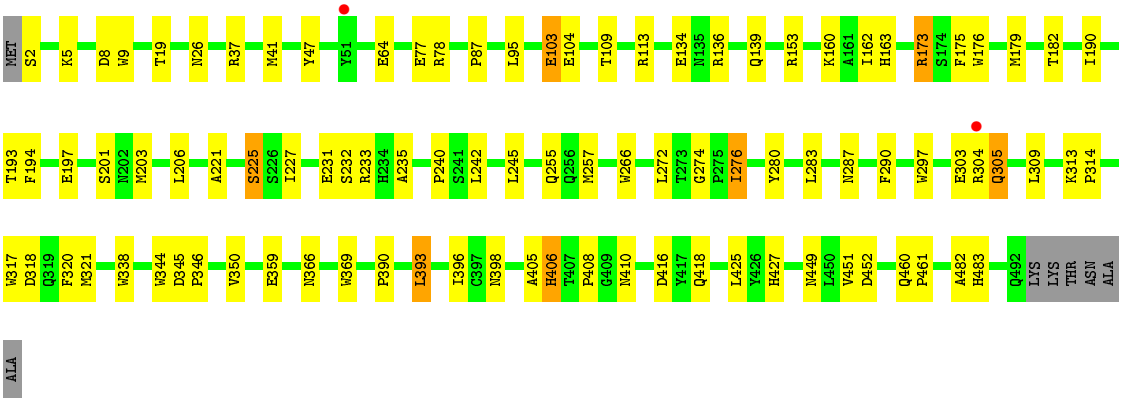
3 Residue-property plots (i)

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

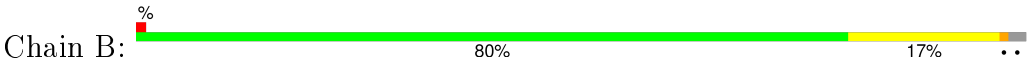
Chain A:

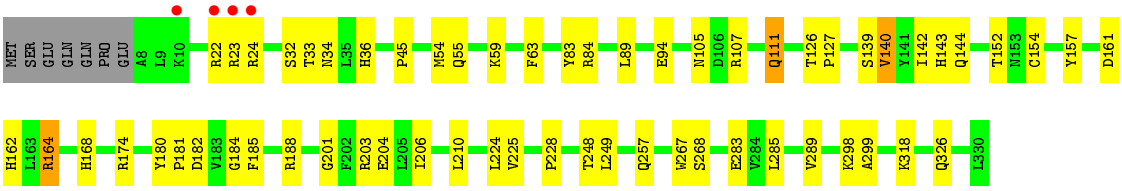




- Molecule 2: Toluene o-xylene monooxygenase component

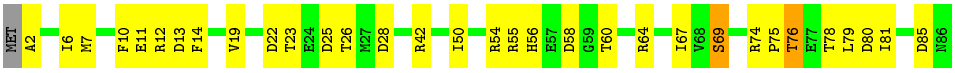
Chain B:





- Molecule 3: Toluene o-xylene monooxygenase component

Chain C:



4 Data and refinement statistics (i)

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
$< I/\sigma(I) >$ [1]	8.17 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.175 , 0.235 0.182 , 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 , 29.9	EDS
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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, 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(^{o})	Ideal(^{o})
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 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	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

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 8.

All (109) 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: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	Interatomic distance (Å)	Clash overlap (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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:GLU:HG2	1:A:231:GLU:OE2	2.15	0.47
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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%)	39 63
2	B	321/330 (97%)	315 (98%)	6 (2%)	0	100 100
3	C	83/86 (96%)	75 (90%)	7 (8%)	1 (1%)	16 31
All	All	894/914 (98%)	850 (95%)	41 (5%)	3 (0%)	46 70

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%)	31	56
2	B	275/282 (98%)	264 (96%)	11 (4%)	38	65
3	C	77/79 (98%)	72 (94%)	5 (6%)	21	40
All	All	769/783 (98%)	733 (95%)	36 (5%)	32	57

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

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Mol	Chain	Res	Type
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 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 (i)

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	0.73	0
6	SO4	A	503	-	4,4,4	0.36	0	6,6,6	0.39	0
6	SO4	B	331	-	4,4,4	0.46	0	6,6,6	0.27	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	502	SO4	1	0
6	B	331	SO4	1	0

5.7 Other polymers (i)

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, 95^{th} 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	2 (0%) 93 92	29, 45, 60, 77	0
2	B	323/330 (97%)	-0.37	4 (1%) 81 78	31, 41, 60, 86	0
3	C	85/86 (98%)	-0.40	0 100 100	40, 53, 65, 74	0
All	All	899/914 (98%)	-0.31	6 (0%) 89 87	29, 45, 61, 86	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	22	ARG	3.5
2	B	23	ARG	3.2
1	A	304	ARG	2.5
2	B	10	LYS	2.5
2	B	24	ARG	2.2
1	A	51	TYR	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

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, 95^{th} 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	LLDF	B-factors(Å^2)	Q<0.9
5	FE	A	501	1/1	0.99	0.17	-0.90	43,43,43,43	0
5	FE	A	500	1/1	1.00	0.17	-0.97	38,38,38,38	0
4	OH	A	499	1/1	1.00	0.16	-1.06	49,49,49,49	0
6	SO4	A	502	5/5	0.92	0.19	-	42,42,46,46	5
6	SO4	B	331	5/5	0.96	0.11	-	47,48,49,49	5
6	SO4	A	503	5/5	0.91	0.16	-	49,51,52,52	5

6.5 Other polymers (i)

There are no such residues in this entry.