Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 07:11 PM GMT

PDB ID	:	1EG9
Title	:	NAPHTHALENE 1,2-DIOXYGENASE WITH INDOLE BOUND IN THE ACTIVE SITE.
Authors	:	Carredano, E.; Karlsson, A.; Kauppi, B.; Choudhury, D.; Parales, R.E.; Parales, J.V.; Lee, K.; Gibson, D.T.; Eklund, H.; Ramaswamy, S.
Deposited on	:	2000-02-15
Resolution	:	1.60 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:

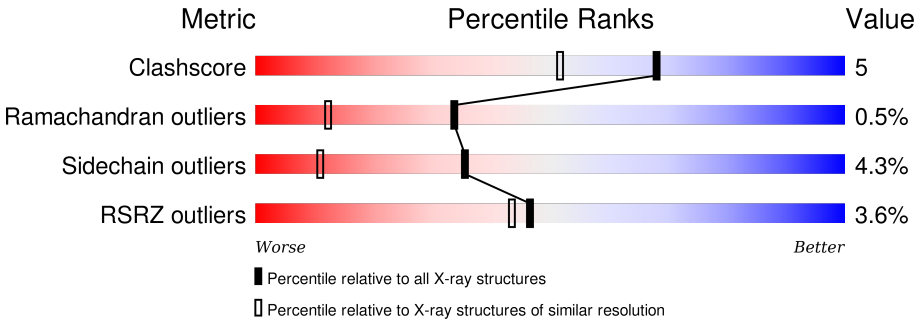
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 1.60 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))
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)
RSRZ outliers	91569	2479 (1.60-1.60)

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	449	2% 86% 12%
2	B	194	6% 76% 18% 5%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	IND	A	708	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5729 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.

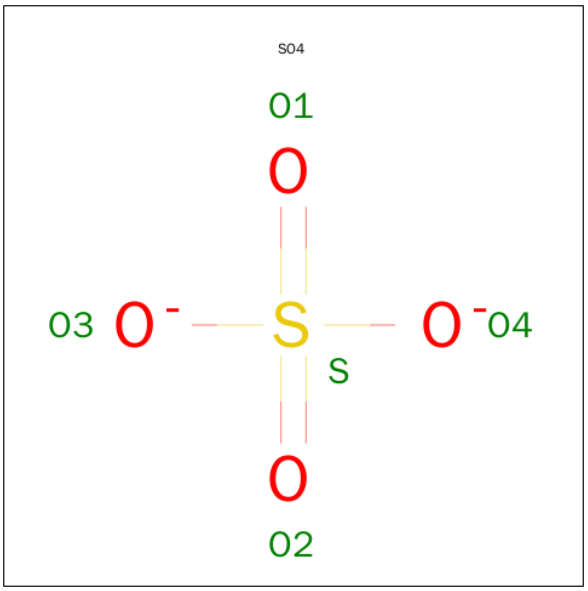
- Molecule 1 is a protein called PROTEIN (NAPHTHALENE 1,2-DIOXYGENASE ALPHA SUBUNIT).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	447	3480	2202	597	665	16	0	0	0

- Molecule 2 is a protein called PROTEIN (NAPHTHALENE 1,2-DIOXYGENASE BETA SUBUNIT).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	193	1608	1007	302	293	6	0	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

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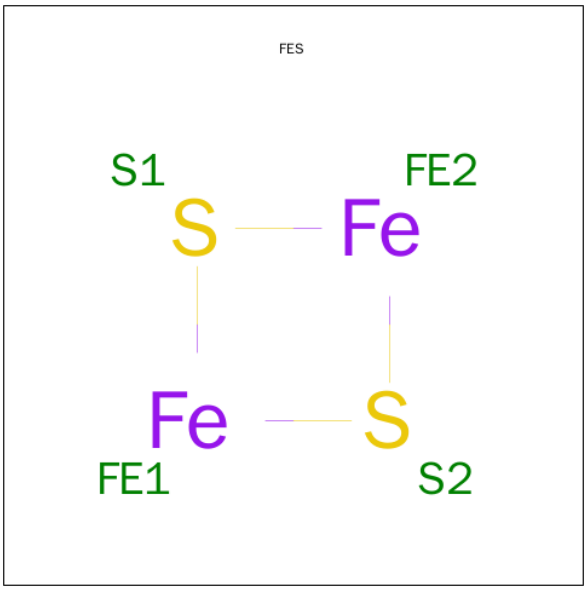
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

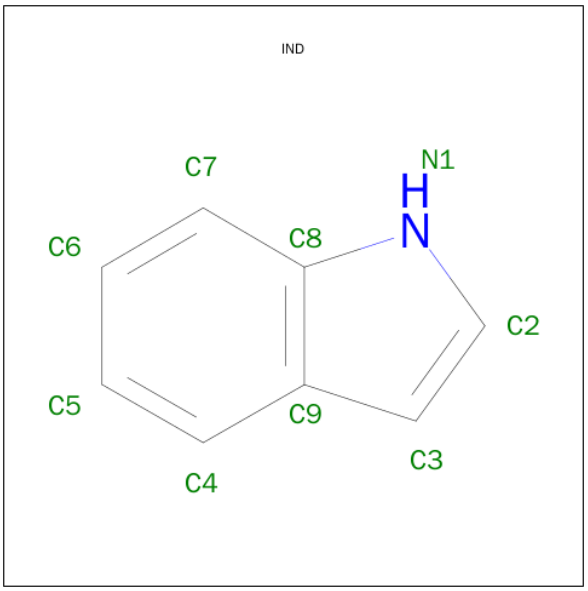
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Fe 1 1	0	0

- Molecule 5 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Fe S 4 2 2	0	0

- Molecule 6 is INDOLE (three-letter code: IND) (formula: C_8H_7N).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C N 9 8 1	0	0

- Molecule 7 is water.

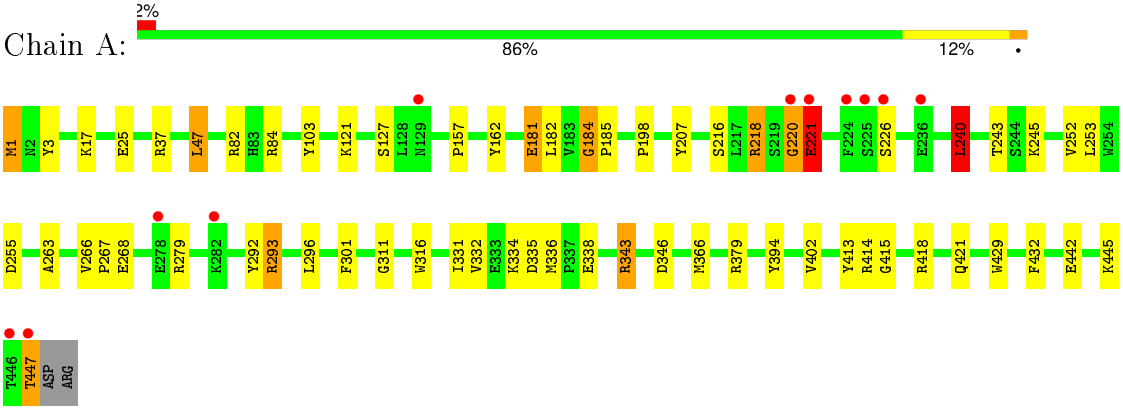
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	430	Total O 430 430	0	0
7	B	182	Total O 182 182	0	0

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.

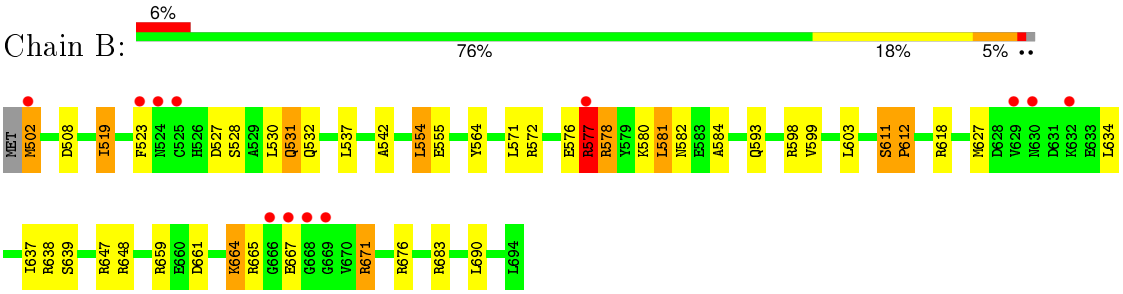
- Molecule 1: PROTEIN (NAPHTHALENE 1,2-DIOXYGENASE ALPHA SUBUNIT)

Chain A:



- Molecule 2: PROTEIN (NAPHTHALENE 1,2-DIOXYGENASE BETA SUBUNIT)

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α, β, γ	140.17Å 140.17Å 209.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 1.60 19.87 – 1.61	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-1.60) 99.4 (19.87-1.61)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.28 (at 1.60Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.190 , 0.220 0.179 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	15.1	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 51.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Outliers	0 of 102099 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5729	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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: IND, SO4, FES, 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	0.63	0/3572	1.39	37/4839 (0.8%)
2	B	0.65	0/1638	1.63	34/2209 (1.5%)
All	All	0.64	0/5210	1.47	71/7048 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	8
2	B	0	4
All	All	0	12

There are no bond length outliers.

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	218	ARG	NE-CZ-NH2	-15.76	112.42	120.30
1	A	218	ARG	NE-CZ-NH1	14.55	127.57	120.30
2	B	618	ARG	NE-CZ-NH1	-14.10	113.25	120.30
2	B	676	ARG	NE-CZ-NH1	13.40	127.00	120.30
2	B	638	ARG	NE-CZ-NH1	10.80	125.70	120.30
2	B	618	ARG	NE-CZ-NH2	10.76	125.68	120.30
2	B	647	ARG	NE-CZ-NH2	10.46	125.53	120.30
1	A	343	ARG	NE-CZ-NH2	-9.86	115.37	120.30
2	B	665	ARG	NE-CZ-NH2	-9.87	115.37	120.30
2	B	659	ARG	NE-CZ-NH2	-9.84	115.38	120.30
1	A	394	TYR	CB-CG-CD2	9.69	126.82	121.00
2	B	638	ARG	NE-CZ-NH2	-9.28	115.66	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	429	TRP	CB-CG-CD1	8.74	138.37	127.00
1	A	218	ARG	CD-NE-CZ	8.25	135.15	123.60
1	A	338	GLU	OE1-CD-OE2	8.06	132.98	123.30
2	B	683	ARG	NE-CZ-NH1	-7.90	116.35	120.30
2	B	572	ARG	NE-CZ-NH2	-7.69	116.46	120.30
2	B	527	ASP	CB-CG-OD1	7.50	125.05	118.30
1	A	184	GLY	CA-C-O	-7.43	107.23	120.60
2	B	598	ARG	NE-CZ-NH1	-7.26	116.67	120.30
1	A	432	PHE	CB-CG-CD1	-7.11	115.82	120.80
1	A	346	ASP	CB-CG-OD1	7.03	124.63	118.30
1	A	185	PRO	CA-N-CD	-6.94	101.78	111.50
2	B	572	ARG	NE-CZ-NH1	6.92	123.76	120.30
2	B	648	ARG	NE-CZ-NH1	-6.87	116.86	120.30
2	B	572	ARG	CD-NE-CZ	6.86	133.20	123.60
2	B	611	SER	CA-C-O	-6.76	105.91	120.10
2	B	612	PRO	N-CA-CB	6.75	111.41	103.30
2	B	671	ARG	NE-CZ-NH2	6.54	123.57	120.30
2	B	612	PRO	CA-N-CD	-6.51	102.38	111.50
1	A	255	ASP	CB-CG-OD1	6.51	124.16	118.30
1	A	394	TYR	CB-CG-CD1	-6.43	117.14	121.00
1	A	207	TYR	CB-CG-CD1	6.39	124.84	121.00
1	A	379	ARG	NE-CZ-NH1	6.37	123.48	120.30
2	B	577	ARG	NE-CZ-NH2	6.28	123.44	120.30
2	B	659	ARG	CD-NE-CZ	-6.27	114.83	123.60
2	B	683	ARG	NE-CZ-NH2	6.20	123.40	120.30
1	A	82	ARG	NE-CZ-NH2	-6.16	117.22	120.30
2	B	564	TYR	CB-CG-CD2	-6.11	117.34	121.00
1	A	292	TYR	CB-CG-CD2	-6.09	117.34	121.00
2	B	647	ARG	NH1-CZ-NH2	-6.08	112.71	119.40
1	A	429	TRP	CD1-CG-CD2	-6.07	101.44	106.30
1	A	279	ARG	NE-CZ-NH2	-5.99	117.31	120.30
1	A	185	PRO	N-CA-CB	5.96	110.45	103.30
2	B	502	MET	CA-CB-CG	5.96	123.43	113.30
1	A	221	GLU	CA-C-O	5.93	132.54	120.10
1	A	413	TYR	CA-CB-CG	5.87	124.56	113.40
1	A	240	LEU	CA-CB-CG	5.80	128.64	115.30
2	B	639	SER	O-C-N	5.77	131.93	122.70
1	A	429	TRP	CE3-CZ3-CH2	-5.68	114.95	121.20
1	A	432	PHE	CB-CG-CD2	5.57	124.70	120.80
2	B	647	ARG	O-C-N	5.51	131.51	122.70
1	A	84	ARG	NE-CZ-NH2	5.48	123.04	120.30
1	A	429	TRP	CD2-CE3-CZ3	5.46	125.91	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	335	ASP	CB-CG-OD1	5.43	123.19	118.30
1	A	293	ARG	CD-NE-CZ	5.42	131.19	123.60
1	A	162	TYR	CA-CB-CG	5.28	123.43	113.40
2	B	612	PRO	N-CD-CG	5.23	111.04	103.20
2	B	578	ARG	CD-NE-CZ	5.19	130.86	123.60
2	B	661	ASP	CB-CG-OD1	-5.15	113.66	118.30
1	A	421	GLN	N-CA-CB	5.15	119.86	110.60
1	A	1	MET	CA-CB-CG	5.12	122.00	113.30
1	A	429	TRP	CB-CG-CD2	-5.12	119.95	126.60
2	B	637	ILE	O-C-N	5.12	130.88	122.70
1	A	157	PRO	N-CA-CB	5.05	109.37	103.30
1	A	293	ARG	NE-CZ-NH1	-5.05	117.77	120.30
2	B	564	TYR	CB-CG-CD1	5.05	124.03	121.00
2	B	676	ARG	NE-CZ-NH2	-5.05	117.78	120.30
1	A	414	ARG	NE-CZ-NH1	-5.04	117.78	120.30
1	A	82	ARG	NE-CZ-NH1	5.04	122.82	120.30
2	B	584	ALA	C-N-CA	-5.01	109.18	121.70

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	127	SER	Mainchain
1	A	184	GLY	Mainchain,Peptide
1	A	220	GLY	Mainchain
1	A	301	PHE	Mainchain,Peptide
1	A	336	MET	Mainchain
1	A	402	VAL	Mainchain
2	B	542	ALA	Mainchain
2	B	580	LYS	Mainchain
2	B	611	SER	Mainchain,Peptide

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	3480	0	3316	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1608	0	1583	22	0
3	A	5	0	0	0	0
3	B	10	0	0	0	0
4	A	1	0	0	0	0
5	A	4	0	0	0	0
6	A	9	0	7	0	0
7	A	430	0	0	7	0
7	B	182	0	0	3	0
All	All	5729	0	4906	47	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 5.

All (47) 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:37:ARG:NH2	7:A:1010:HOH:O	2.05	0.81
1:A:181:GLU:HG2	1:A:334:LYS:HG2	1.62	0.81
1:A:240:LEU:HD13	1:A:252:VAL:HB	1.68	0.74
2:B:576:GLU:HA	2:B:577:ARG:NH1	2.02	0.73
2:B:519:ILE:HD13	2:B:523:PHE:CZ	2.27	0.69
2:B:577:ARG:HH21	2:B:578:ARG:HE	1.40	0.68
1:A:266:VAL:HB	1:A:267:PRO:HD3	1.74	0.67
1:A:220:GLY:HA3	7:A:894:HOH:O	1.98	0.62
2:B:577:ARG:NH2	2:B:578:ARG:HE	2.02	0.57
2:B:577:ARG:HH21	2:B:578:ARG:NE	2.01	0.56
2:B:634:LEU:HD23	2:B:664:LYS:HB2	1.88	0.56
1:A:245:LYS:O	1:A:418:ARG:NH2	2.40	0.55
1:A:47:LEU:HD22	1:A:182:LEU:HD23	1.88	0.55
2:B:531:GLN:NE2	2:B:532:GLN:OE1	2.41	0.53
1:A:181:GLU:HG3	1:A:332:VAL:HG23	1.89	0.53
2:B:577:ARG:HD2	2:B:578:ARG:HG3	1.90	0.53
2:B:576:GLU:HA	2:B:577:ARG:HH11	1.72	0.52
1:A:181:GLU:CG	1:A:334:LYS:HG2	2.37	0.51
1:A:366:MET:HE2	7:A:1180:HOH:O	2.11	0.49
2:B:555:GLU:HG3	7:B:843:HOH:O	2.11	0.49
2:B:519:ILE:HG23	2:B:523:PHE:CE1	2.48	0.49
1:A:447:THR:HB	7:A:1034:HOH:O	2.12	0.49
1:A:221:GLU:O	1:A:263:ALA:N	2.46	0.48
1:A:216:SER:O	1:A:220:GLY:HA3	2.13	0.48
2:B:530:LEU:HD21	2:B:627:MET:HG2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:ARG:HD3	7:A:1149:HOH:O	2.15	0.47
2:B:577:ARG:HH21	2:B:578:ARG:HB2	1.80	0.47
2:B:554:LEU:HD22	2:B:599:VAL:HG21	1.99	0.45
2:B:519:ILE:HD13	2:B:523:PHE:CE1	2.50	0.45
1:A:311:GLY:HA2	1:A:331:ILE:HG13	1.96	0.45
2:B:577:ARG:NH2	2:B:578:ARG:NE	2.62	0.45
1:A:25:GLU:HB3	7:A:1079:HOH:O	2.16	0.45
1:A:252:VAL:HG22	1:A:296:LEU:HD22	1.99	0.45
2:B:577:ARG:H	2:B:577:ARG:CZ	2.30	0.45
2:B:528:SER:O	2:B:532:GLN:NE2	2.51	0.44
1:A:216:SER:O	1:A:220:GLY:N	2.51	0.43
2:B:577:ARG:N	2:B:577:ARG:NE	2.67	0.42
1:A:198:PRO:HB2	1:A:316:TRP:CE2	2.55	0.42
1:A:442:GLU:O	1:A:445:LYS:HG2	2.20	0.42
2:B:537:LEU:HD12	7:B:750:HOH:O	2.18	0.42
1:A:243:THR:OG1	1:A:415:GLY:HA3	2.20	0.42
1:A:1:MET:HG2	1:A:3:TYR:CE2	2.55	0.41
1:A:37:ARG:HB2	7:A:1007:HOH:O	2.20	0.41
2:B:581:LEU:O	2:B:582:ASN:C	2.59	0.41
1:A:17:LYS:HD3	1:A:17:LYS:HA	1.87	0.41
2:B:664:LYS:HG2	7:B:880:HOH:O	2.21	0.40
1:A:268:GLU:CD	1:A:343:ARG:HH22	2.25	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	445/449 (99%)	429 (96%)	14 (3%)	2 (0%)	39	17
2	B	191/194 (98%)	184 (96%)	6 (3%)	1 (0%)	34	12
All	All	636/643 (99%)	613 (96%)	20 (3%)	3 (0%)	34	12

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	221	GLU
1	A	226	SER
2	B	612	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	367/369 (100%)	358 (98%)	9 (2%)	55 26
2	B	172/173 (99%)	158 (92%)	14 (8%)	15 2
All	All	539/542 (99%)	516 (96%)	23 (4%)	35 11

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	LEU
1	A	103	TYR
1	A	121	LYS
1	A	181	GLU
1	A	221	GLU
1	A	240	LEU
1	A	253	LEU
1	A	293	ARG
1	A	447	THR
2	B	502	MET
2	B	508	ASP
2	B	519	ILE
2	B	531	GLN
2	B	554	LEU
2	B	571	LEU
2	B	577	ARG
2	B	581	LEU
2	B	593	GLN
2	B	603	LEU
2	B	664	LYS

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Mol	Chain	Res	Type
2	B	667	GLU
2	B	671	ARG
2	B	690	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	115	GLN
2	B	514	HIS
2	B	531	GLN
2	B	532	GLN
2	B	556	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

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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	IND	A	708	-	8,10,10	0.76	0	9,13,13	1.26	1 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	SO4	A	717	-	4,4,4	3.39	2 (50%)	6,6,6	0.95	0
5	FES	A	751	1	0,4,4	0.00	-	0,4,4	0.00	-
3	SO4	B	702	-	4,4,4	0.92	0	6,6,6	0.10	0
3	SO4	B	707	-	4,4,4	0.91	0	6,6,6	0.40	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	IND	A	708	-	-	0/0/0/0	0/2/2/2
3	SO4	A	717	-	-	0/0/0/0	0/0/0/0
5	FES	A	751	1	-	0/0/4/4	0/1/1/1
3	SO4	B	702	-	-	0/0/0/0	0/0/0/0
3	SO4	B	707	-	-	0/0/0/0	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	717	SO4	O3-S	-4.84	1.30	1.47
3	A	717	SO4	O1-S	4.65	1.63	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	A	708	IND	C6-C7-C8	-2.38	116.25	120.06

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

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	447/449 (99%)	-0.32	11 (2%)	61	58	11, 16, 35, 55	0
2	B	193/194 (99%)	-0.14	12 (6%)	24	21	11, 16, 40, 56	0
All	All	640/643 (99%)	-0.27	23 (3%)	46	43	11, 16, 37, 56	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	447	THR	12.8
1	A	446	THR	9.0
1	A	236	GLU	5.9
2	B	525	CYS	5.4
2	B	523	PHE	5.2
2	B	577	ARG	4.3
2	B	502	MET	4.2
1	A	221	GLU	4.1
2	B	669	GLY	3.9
2	B	524	ASN	3.5
2	B	629	VAL	3.2
2	B	666	GLY	3.2
1	A	226	SER	3.0
1	A	129	ASN	2.7
2	B	630	ASN	2.5
1	A	282	LYS	2.4
1	A	225	SER	2.2
2	B	667	GLU	2.2
1	A	224	PHE	2.1
1	A	278	GLU	2.1
2	B	632	LYS	2.1
2	B	668	GLY	2.1
1	A	220	GLY	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 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. 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
6	IND	A	708	9/9	0.81	0.15	3.25	32,34,36,36	0
3	SO4	B	707	5/5	0.94	0.12	0.08	39,40,41,42	0
5	FES	A	751	4/4	0.99	0.05	-1.12	8,10,16,17	0
4	FE	A	752	1/1	0.99	0.09	-	22,22,22,22	0
3	SO4	A	717	5/5	0.96	0.39	-	20,20,20,20	0
3	SO4	B	702	5/5	0.81	0.21	-	70,70,71,71	0

6.5 Other polymers (i)

There are no such residues in this entry.