



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:02 AM GMT

PDB ID : 2VQB
Title : BACTERIAL FLAVIN-CONTAINING MONOOXYGENASE IN COMPLEX
WITH NADP: SOAKING IN AERATED SOLUTION
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Deposited on : 2008-03-12
Resolution : 2.80 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

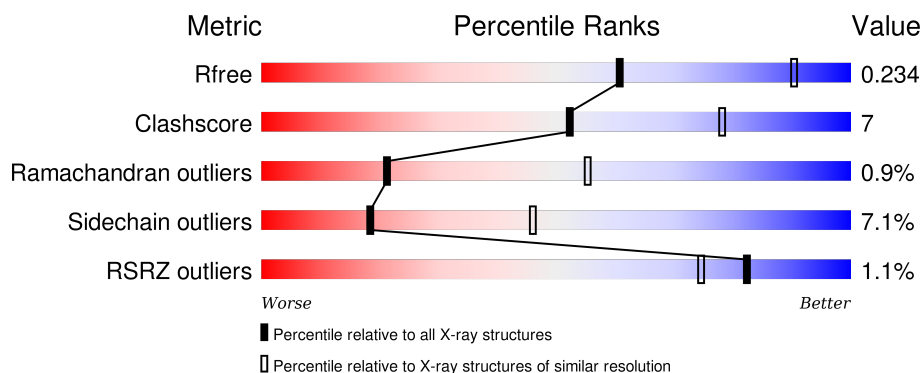
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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 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 $\leq 5\%$. 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	461	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>16%</div> <div>• •</div> </div> </div>
1	B	461	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>• •</div> </div> </div>
1	C	461	<div> <div>%</div> <div> <div></div> <div>79%</div> <div>16%</div> <div>• •</div> </div> </div>
1	D	461	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>• •</div> </div> </div>

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
5	EPE	A	1452	-	-	-	X
5	EPE	B	1452	-	-	-	X
5	EPE	C	1451	-	-	-	X
5	EPE	D	1452	-	-	-	X
6	OXY	D	1454	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 15036 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.

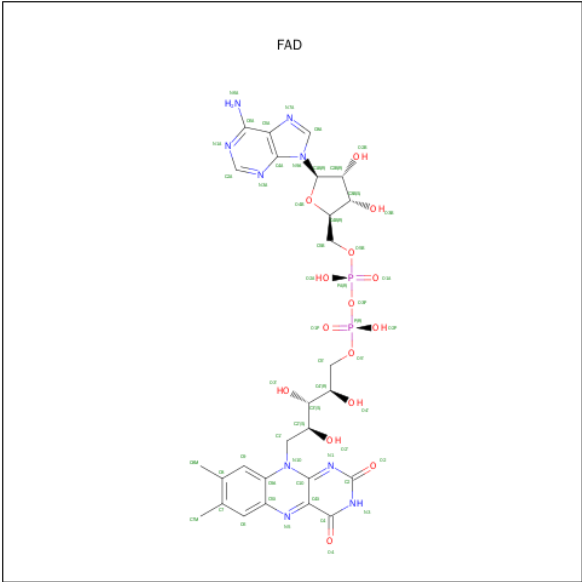
- Molecule 1 is a protein called FLAVIN-CONTAINING MONOOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	443	Total	C	N	O	S	0	0	0
			3638	2335	603	680	20			
1	B	443	Total	C	N	O	S	0	0	0
			3638	2335	603	680	20			
1	C	443	Total	C	N	O	S	0	0	0
			3638	2335	603	680	20			
1	D	443	Total	C	N	O	S	0	0	0
			3638	2335	603	680	20			

There are 8 discrepancies between the modelled and reference sequences:

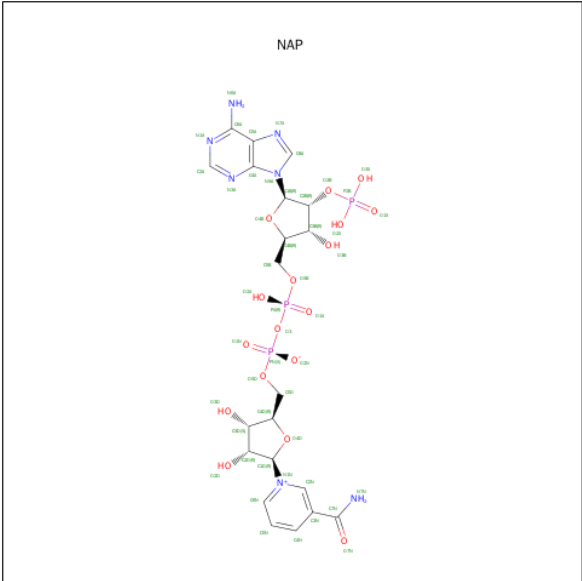
Chain	Residue	Modelled	Actual	Comment	Reference
A	158	ALA	GLU	ENGINEERED MUTATION	UNP Q83XK4
A	159	ALA	GLU	ENGINEERED MUTATION	UNP Q83XK4
B	158	ALA	GLU	ENGINEERED MUTATION	UNP Q83XK4
B	159	ALA	GLU	ENGINEERED MUTATION	UNP Q83XK4
C	158	ALA	GLU	ENGINEERED MUTATION	UNP Q83XK4
C	159	ALA	GLU	ENGINEERED MUTATION	UNP Q83XK4
D	158	ALA	GLU	ENGINEERED MUTATION	UNP Q83XK4
D	159	ALA	GLU	ENGINEERED MUTATION	UNP Q83XK4

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).

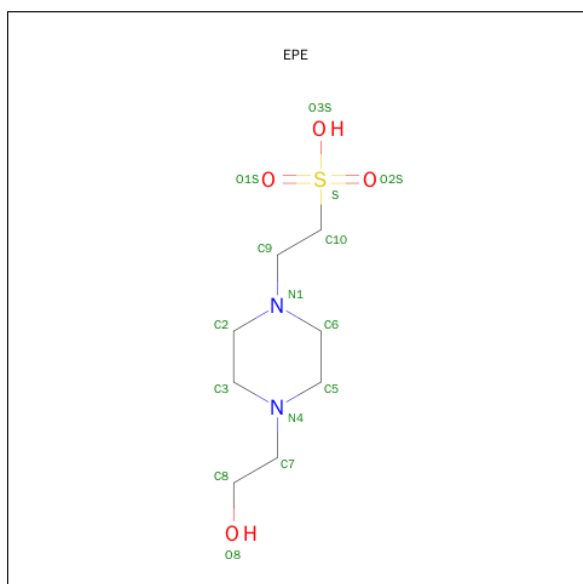


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		
4	A	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		
4	C	1	Total	Cl	0	0
			1	1		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



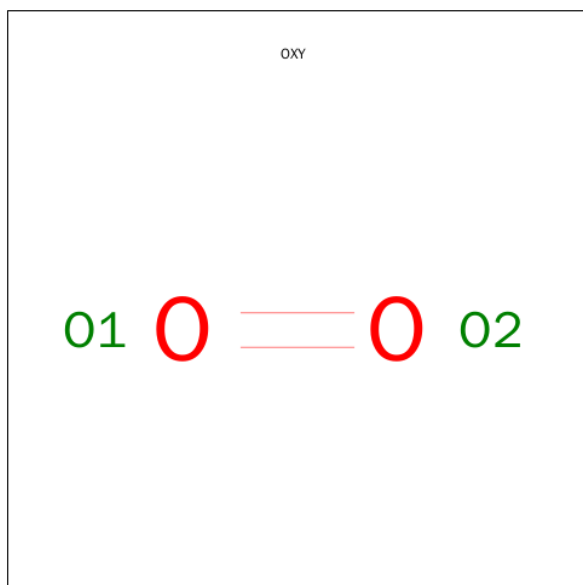
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
5	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
5	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 6 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O₂).

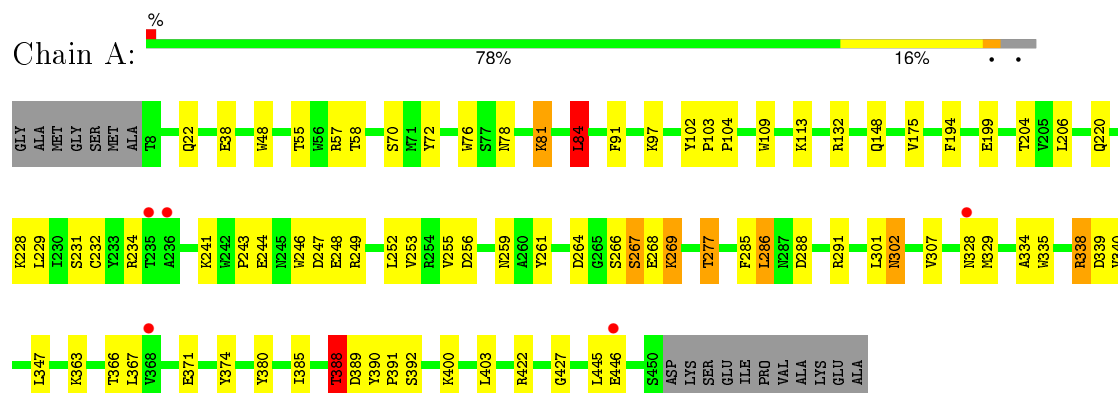


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	O	0	0
			2	2		
6	A	1	Total	O	0	0
			2	2		
6	B	1	Total	O	0	0
			2	2		
6	B	1	Total	O	0	0
			2	2		
6	C	1	Total	O	0	0
			2	2		
6	C	1	Total	O	0	0
			2	2		
6	D	1	Total	O	0	0
			2	2		
6	D	1	Total	O	0	0
			2	2		

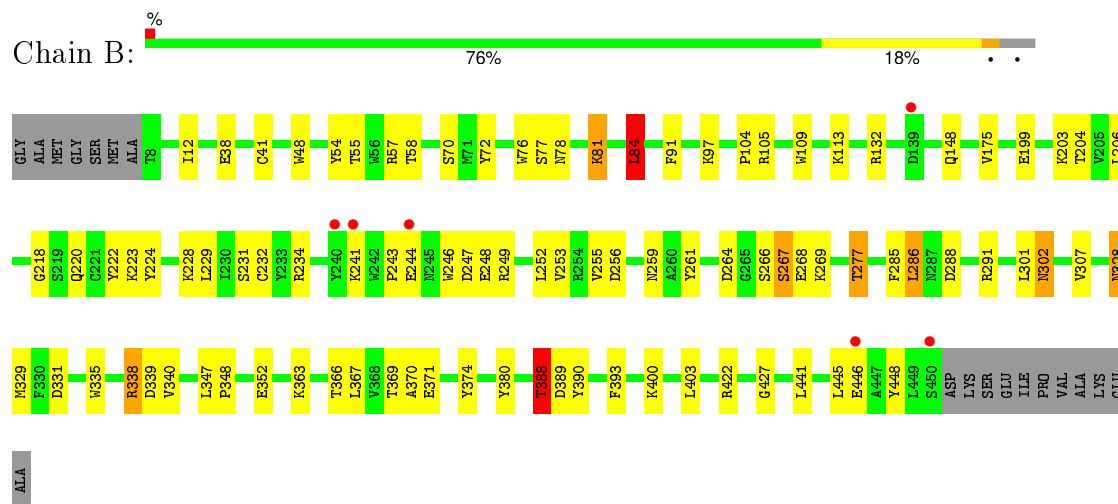
3 Residue-property plots

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.

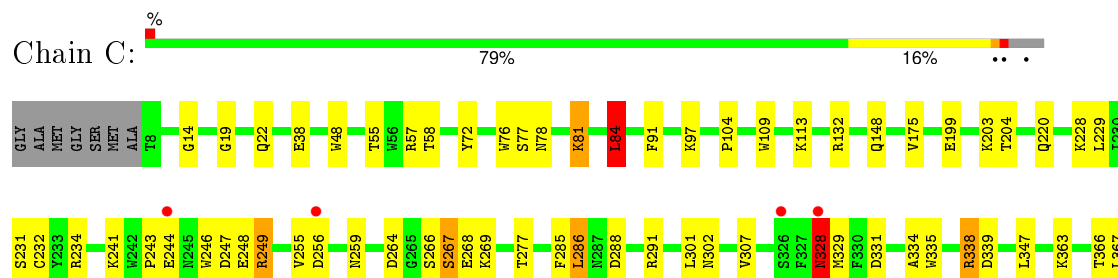
• Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE

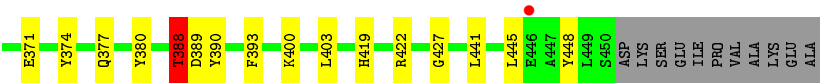


• Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE

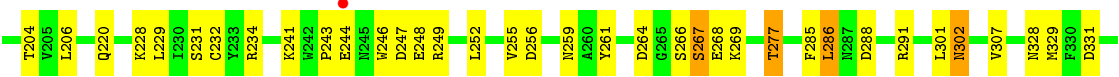
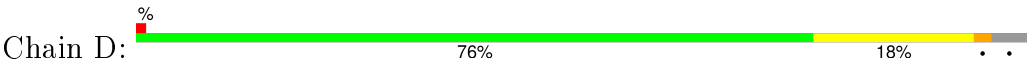


• Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE





● Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	218.66Å 218.66Å 130.68Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.72 – 2.80 29.72 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.72-2.80) 99.6 (29.72-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.226 , 0.242 0.219 , 0.234	Depositor DCC
R_{free} test set	4332 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	61.7	Xtriage
Anisotropy	0.317	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 21.1	EDS
Estimated twinning fraction	0.019 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 86900 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15036	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 2.34% 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, 0.375 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: NAP, EPE, OXY, FAD, CL

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	0.61	0/3752	0.67	2/5095 (0.0%)
1	B	0.63	0/3752	0.68	2/5095 (0.0%)
1	C	0.61	0/3752	0.68	2/5095 (0.0%)
1	D	0.59	0/3752	0.66	2/5095 (0.0%)
All	All	0.61	0/15008	0.67	8/20380 (0.0%)

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	1
1	B	0	1
1	C	0	1
1	D	0	1
All	All	0	4

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	84	LEU	CA-CB-CG	7.13	131.71	115.30
1	C	445	LEU	CA-CB-CG	6.65	130.59	115.30
1	D	445	LEU	CA-CB-CG	6.61	130.50	115.30
1	D	84	LEU	CA-CB-CG	6.54	130.34	115.30
1	C	84	LEU	CA-CB-CG	6.41	130.05	115.30
1	B	84	LEU	CA-CB-CG	6.38	129.98	115.30
1	A	445	LEU	CA-CB-CG	6.34	129.88	115.30
1	B	445	LEU	CA-CB-CG	6.11	129.34	115.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	388	THR	Peptide
1	B	388	THR	Peptide
1	C	388	THR	Peptide
1	D	388	THR	Peptide

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	3638	0	3428	46	0
1	B	3638	0	3428	55	0
1	C	3638	0	3428	47	0
1	D	3638	0	3428	49	0
2	A	53	0	31	1	0
2	B	53	0	31	1	0
2	C	53	0	31	0	0
2	D	53	0	31	1	0
3	A	48	0	25	2	0
3	B	48	0	25	3	0
3	C	48	0	25	2	0
3	D	48	0	25	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	15	0	17	4	0
5	B	15	0	17	5	0
5	C	15	0	17	0	0
5	D	15	0	17	0	0
6	A	4	0	0	0	0
6	B	4	0	0	0	0
6	C	4	0	0	0	0
6	D	4	0	0	0	0
All	All	15036	0	14004	195	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 7.

All (195) 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)
5:A:1452:EPE:H21	5:B:1452:EPE:H21	1.21	1.17
1:C:388:THR:HG22	1:C:390:TYR:H	1.36	0.89
1:C:422:ARG:HD3	1:C:427:GLY:O	1.79	0.82
1:D:422:ARG:HD3	1:D:427:GLY:O	1.80	0.81
1:A:335:TRP:HB3	1:A:388:THR:HG21	1.63	0.81
1:A:388:THR:HG22	1:A:390:TYR:H	1.46	0.80
1:B:422:ARG:HD3	1:B:427:GLY:O	1.81	0.79
1:A:338:ARG:NH1	1:A:339:ASP:OD1	2.15	0.79
1:D:388:THR:HG22	1:D:390:TYR:H	1.48	0.78
1:B:335:TRP:HB3	1:B:388:THR:HG21	1.65	0.77
1:C:335:TRP:HB3	1:C:388:THR:HG21	1.66	0.76
1:B:388:THR:HG22	1:B:390:TYR:H	1.51	0.75
1:A:285:PHE:CD1	1:A:286:LEU:HD13	2.21	0.75
1:D:335:TRP:HB3	1:D:388:THR:HG21	1.69	0.74
1:B:57:ARG:NH2	1:C:175:VAL:O	2.20	0.74
1:D:285:PHE:CD1	1:D:286:LEU:HD13	2.23	0.73
1:A:422:ARG:HD3	1:A:427:GLY:O	1.88	0.72
1:D:338:ARG:NH1	1:D:339:ASP:OD1	2.23	0.71
1:B:338:ARG:NH1	1:B:339:ASP:OD1	2.23	0.71
1:A:57:ARG:NH2	1:D:175:VAL:O	2.24	0.71
1:A:301:LEU:O	1:A:302:ASN:HB2	1.89	0.70
1:B:329:MET:HE2	1:B:380:TYR:HE2	1.55	0.70
1:B:175:VAL:O	1:C:57:ARG:NH2	2.25	0.70
1:C:329:MET:HE2	1:C:380:TYR:HE2	1.57	0.69
1:B:285:PHE:CD1	1:B:286:LEU:HD13	2.27	0.69
1:D:329:MET:HE2	1:D:380:TYR:HE2	1.58	0.68
1:C:285:PHE:CD1	1:C:286:LEU:HD13	2.28	0.68
1:C:338:ARG:NH1	1:C:339:ASP:OD1	2.28	0.67
1:D:301:LEU:O	1:D:302:ASN:HB2	1.96	0.65
1:C:301:LEU:O	1:C:302:ASN:HB2	1.96	0.65
1:A:329:MET:HE2	1:A:380:TYR:HE2	1.62	0.65
1:C:388:THR:CG2	1:C:390:TYR:H	2.09	0.64
1:A:175:VAL:O	1:D:57:ARG:NH2	2.31	0.63
1:B:264:ASP:HB3	1:B:266:SER:H	1.63	0.63
1:B:301:LEU:O	1:B:302:ASN:HB2	1.97	0.63
1:A:264:ASP:HB3	1:A:266:SER:H	1.64	0.62
1:C:264:ASP:HB3	1:C:266:SER:H	1.65	0.62
1:B:256:ASP:HB3	1:B:259:ASN:H	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:422:ARG:CD	1:B:427:GLY:O	2.47	0.61
1:D:264:ASP:HB3	1:D:266:SER:H	1.65	0.61
1:C:422:ARG:CD	1:C:427:GLY:O	2.48	0.61
1:D:256:ASP:HB3	1:D:259:ASN:H	1.67	0.60
1:C:81:LYS:HG2	1:C:91:PHE:CE2	2.37	0.60
1:D:204:THR:HG23	1:D:228:LYS:HB3	1.83	0.60
1:A:204:THR:HG23	1:A:228:LYS:HB3	1.85	0.59
5:A:1452:EPE:H102	5:B:1452:EPE:H62	1.84	0.59
1:C:256:ASP:HB3	1:C:259:ASN:H	1.67	0.59
1:B:329:MET:HE2	1:B:380:TYR:CE2	2.37	0.59
1:A:256:ASP:HB3	1:A:259:ASN:H	1.66	0.59
1:D:388:THR:CG2	1:D:390:TYR:H	2.14	0.59
1:A:388:THR:CG2	1:A:390:TYR:H	2.14	0.58
1:C:363:LYS:O	1:C:366:THR:HB	2.03	0.58
1:B:204:THR:HG23	1:B:228:LYS:HB3	1.85	0.58
1:D:81:LYS:HG2	1:D:91:PHE:CE2	2.39	0.58
1:D:422:ARG:CD	1:D:427:GLY:O	2.50	0.58
1:B:81:LYS:HG2	1:B:91:PHE:CE2	2.39	0.57
1:D:363:LYS:O	1:D:366:THR:HB	2.03	0.57
1:B:277:THR:HG22	3:B:501:NAP:C4A	2.34	0.57
1:A:363:LYS:O	1:A:366:THR:HB	2.05	0.57
1:B:363:LYS:O	1:B:366:THR:HB	2.04	0.57
1:C:277:THR:HG22	3:C:501:NAP:C4A	2.35	0.56
1:A:76:TRP:HA	1:A:104:PRO:HA	1.87	0.56
1:B:76:TRP:HA	1:B:104:PRO:HA	1.88	0.56
5:A:1452:EPE:H21	5:B:1452:EPE:C2	2.15	0.56
1:B:338:ARG:HG3	1:B:339:ASP:N	2.21	0.56
1:B:388:THR:CG2	1:B:390:TYR:H	2.17	0.56
1:D:285:PHE:HD1	1:D:286:LEU:HD13	1.70	0.55
1:C:329:MET:HE2	1:C:380:TYR:CE2	2.38	0.55
1:D:329:MET:HE2	1:D:380:TYR:CE2	2.38	0.55
1:D:277:THR:HG22	3:D:501:NAP:C4A	2.37	0.55
1:A:55:THR:OG1	1:A:57:ARG:HG3	2.07	0.54
1:A:232:CYS:HA	1:A:249:ARG:O	2.07	0.54
1:D:76:TRP:HA	1:D:104:PRO:HA	1.90	0.54
1:C:109:TRP:CE2	1:C:113:LYS:HD2	2.43	0.54
1:A:329:MET:HE2	1:A:380:TYR:CE2	2.42	0.53
1:A:422:ARG:CD	1:A:427:GLY:O	2.54	0.53
1:C:204:THR:HG23	1:C:228:LYS:HB3	1.91	0.53
1:C:76:TRP:HA	1:C:104:PRO:HA	1.90	0.53
1:A:285:PHE:HD1	1:A:286:LEU:HD13	1.71	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:TRP:CE2	1:B:113:LYS:HD2	2.44	0.52
1:D:109:TRP:CE2	1:D:113:LYS:HD2	2.45	0.52
1:A:81:LYS:HG2	1:A:91:PHE:CE2	2.45	0.51
1:D:72:TYR:OH	3:D:501:NAP:H4N	2.10	0.51
1:B:132:ARG:HH21	1:B:148:GLN:HE21	1.58	0.51
1:A:277:THR:HG22	3:A:501:NAP:C4A	2.40	0.51
1:A:229:LEU:O	1:A:246:TRP:HA	2.10	0.51
1:B:253:VAL:HG22	1:B:261:TYR:O	2.11	0.51
1:A:109:TRP:CE2	1:A:113:LYS:HD2	2.46	0.51
1:D:338:ARG:HG3	1:D:339:ASP:N	2.26	0.51
1:C:267:SER:O	1:C:268:GLU:HG2	2.11	0.50
1:C:371:GLU:HA	1:C:374:TYR:CZ	2.47	0.50
1:B:132:ARG:HH21	1:B:148:GLN:NE2	2.09	0.50
1:A:338:ARG:HG3	1:A:339:ASP:N	2.27	0.50
1:B:288:ASP:OD1	1:B:291:ARG:NH1	2.44	0.50
1:C:132:ARG:HH21	1:C:148:GLN:NE2	2.10	0.49
1:D:288:ASP:OD1	1:D:291:ARG:NH1	2.45	0.49
1:C:72:TYR:OH	3:C:501:NAP:H4N	2.12	0.49
5:A:1452:EPE:C2	5:B:1452:EPE:H21	2.15	0.49
1:C:232:CYS:HA	1:C:249:ARG:O	2.13	0.49
1:A:288:ASP:OD1	1:A:291:ARG:NH1	2.46	0.49
1:C:288:ASP:OD1	1:C:291:ARG:NH1	2.46	0.49
1:B:81:LYS:O	1:B:84:LEU:HD22	2.13	0.48
1:B:231:SER:HB2	1:B:248:GLU:HG2	1.94	0.48
1:D:256:ASP:HB2	1:D:261:TYR:HE1	1.76	0.48
1:D:371:GLU:HA	1:D:374:TYR:CZ	2.49	0.48
1:D:206:LEU:HD21	1:D:252:LEU:HD22	1.96	0.48
1:C:338:ARG:HG3	1:C:339:ASP:N	2.29	0.48
1:D:81:LYS:O	1:D:84:LEU:HD22	2.14	0.48
1:C:388:THR:HG22	1:C:390:TYR:N	2.18	0.48
1:B:285:PHE:HD1	1:B:286:LEU:HD13	1.76	0.48
1:B:203:LYS:HZ2	1:C:203:LYS:HZ1	1.61	0.48
1:D:14:GLY:O	1:D:19:GLY:HA3	2.14	0.47
1:B:371:GLU:HA	1:B:374:TYR:CZ	2.49	0.47
1:A:72:TYR:OH	3:A:501:NAP:H4N	2.15	0.47
1:B:55:THR:OG1	1:B:57:ARG:HG3	2.14	0.47
1:D:232:CYS:HA	1:D:249:ARG:O	2.14	0.47
1:C:109:TRP:NE1	1:C:113:LYS:HD2	2.29	0.47
1:A:253:VAL:HG22	1:A:261:TYR:O	2.15	0.47
1:C:285:PHE:HD1	1:C:286:LEU:HD13	1.79	0.47
1:D:55:THR:OG1	1:D:57:ARG:HG3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:256:ASP:HB2	1:B:261:TYR:HE1	1.79	0.46
1:A:72:TYR:HB3	1:A:194:PHE:O	2.15	0.46
1:C:81:LYS:O	1:C:84:LEU:HD22	2.15	0.46
1:D:231:SER:HB2	1:D:248:GLU:HG2	1.96	0.46
1:B:232:CYS:HA	1:B:249:ARG:O	2.15	0.46
1:B:331:ASP:HB3	1:B:390:TYR:CE1	2.51	0.46
1:B:206:LEU:HD21	1:B:252:LEU:HD22	1.97	0.46
1:B:104:PRO:HD3	1:B:448:TYR:CE1	2.51	0.46
1:D:229:LEU:O	1:D:246:TRP:HA	2.15	0.46
1:C:335:TRP:HD1	1:C:388:THR:HG23	1.81	0.46
1:B:277:THR:HG22	3:B:501:NAP:C5A	2.46	0.45
1:A:335:TRP:HB3	1:A:388:THR:CG2	2.42	0.45
5:B:1452:EPE:H81	5:B:1452:EPE:H31	1.85	0.45
1:A:256:ASP:HB2	1:A:261:TYR:HE1	1.81	0.45
1:D:132:ARG:HH21	1:D:148:GLN:HE21	1.65	0.45
1:B:335:TRP:HB3	1:B:388:THR:CG2	2.41	0.45
1:B:223:LYS:HE2	1:B:224:TYR:CE1	2.52	0.45
1:C:231:SER:HB2	1:C:248:GLU:HG2	1.98	0.44
1:C:104:PRO:HD3	1:C:448:TYR:CE1	2.51	0.44
1:B:328:ASN:OD1	1:B:393:PHE:CZ	2.71	0.44
1:A:206:LEU:HD21	1:A:252:LEU:HD22	1.99	0.44
1:B:109:TRP:NE1	1:B:113:LYS:HD2	2.32	0.44
1:A:132:ARG:HH21	1:A:148:GLN:NE2	2.16	0.44
1:A:335:TRP:HD1	1:A:388:THR:HG23	1.83	0.44
1:B:267:SER:O	1:B:268:GLU:HG2	2.17	0.44
1:D:109:TRP:NE1	1:D:113:LYS:HD2	2.33	0.43
1:B:229:LEU:O	1:B:246:TRP:HA	2.18	0.43
1:C:76:TRP:O	1:C:77:SER:C	2.56	0.43
1:B:203:LYS:NZ	1:C:203:LYS:NZ	2.66	0.43
1:B:70:SER:HB2	2:B:500:FAD:HM82	2.00	0.43
1:D:267:SER:O	1:D:268:GLU:HG2	2.18	0.43
1:B:203:LYS:NZ	1:C:203:LYS:HZ1	2.15	0.43
1:A:371:GLU:HA	1:A:374:TYR:CZ	2.54	0.43
1:B:72:TYR:OH	3:B:501:NAP:H4N	2.19	0.43
1:D:132:ARG:HH21	1:D:148:GLN:NE2	2.16	0.43
1:D:9:ARG:HG3	1:D:38:GLU:HB2	2.01	0.43
1:A:301:LEU:O	1:A:302:ASN:CB	2.62	0.43
1:A:81:LYS:O	1:A:84:LEU:HD22	2.18	0.43
1:C:229:LEU:O	1:C:246:TRP:HA	2.19	0.42
1:B:222:TYR:HB2	1:B:229:LEU:HD11	2.02	0.42
1:D:331:ASP:HB3	1:D:390:TYR:CE1	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:TRP:O	1:B:77:SER:C	2.56	0.42
1:B:218:GLY:HA3	1:B:246:TRP:CH2	2.55	0.42
1:B:369:THR:O	1:B:370:ALA:C	2.56	0.42
1:B:335:TRP:HD1	1:B:388:THR:HG23	1.85	0.42
1:D:348:PRO:HB2	1:D:352:GLU:HB2	2.01	0.42
1:A:267:SER:O	1:A:268:GLU:HG2	2.19	0.42
1:C:14:GLY:O	1:C:19:GLY:HA3	2.19	0.42
1:C:331:ASP:HB3	1:C:390:TYR:CE1	2.55	0.42
1:B:54:TYR:HB2	1:B:105:ARG:CZ	2.50	0.42
1:C:419:HIS:N	1:C:419:HIS:CD2	2.88	0.42
1:D:369:THR:O	1:D:370:ALA:C	2.58	0.42
1:A:385:ILE:HD12	1:A:392:SER:HA	2.01	0.42
1:A:70:SER:HB2	2:A:500:FAD:HM82	2.02	0.41
1:A:132:ARG:HH21	1:A:148:GLN:HE21	1.68	0.41
1:C:328:ASN:OD1	1:C:393:PHE:CZ	2.73	0.41
1:B:12:ILE:O	1:B:41:CYS:HA	2.20	0.41
1:D:70:SER:HB2	2:D:500:FAD:HM82	2.03	0.41
1:D:76:TRP:O	1:D:77:SER:C	2.57	0.41
1:B:348:PRO:HB2	1:B:352:GLU:HB2	2.03	0.41
1:D:335:TRP:HD1	1:D:388:THR:HG23	1.86	0.41
1:A:22:GLN:HA	1:A:334:ALA:HB1	2.02	0.41
1:D:335:TRP:HB3	1:D:388:THR:CG2	2.44	0.41
1:C:55:THR:OG1	1:C:57:ARG:HG3	2.21	0.41
1:C:132:ARG:HH21	1:C:148:GLN:HE21	1.67	0.41
1:C:329:MET:HE3	1:C:377:GLN:HG3	2.03	0.41
1:A:269:LYS:HE3	1:A:269:LYS:HB2	1.87	0.41
1:D:97:LYS:NZ	1:D:449:LEU:O	2.54	0.41
1:A:390:TYR:HA	1:A:391:PRO:HD3	1.96	0.41
1:A:102:TYR:N	1:A:103:PRO:HD3	2.36	0.40
1:D:12:ILE:O	1:D:41:CYS:HA	2.21	0.40
1:C:22:GLN:HA	1:C:334:ALA:HB1	2.03	0.40
1:D:22:GLN:HA	1:D:334:ALA:HB1	2.03	0.40
1:D:329:MET:HE3	1:D:377:GLN:HG3	2.03	0.40
1:A:231:SER:HB2	1:A:248:GLU:HG2	2.03	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	441/461 (96%)	410 (93%)	27 (6%)	4 (1%)	21	55
1	B	441/461 (96%)	408 (92%)	29 (7%)	4 (1%)	21	55
1	C	441/461 (96%)	408 (92%)	29 (7%)	4 (1%)	21	55
1	D	441/461 (96%)	407 (92%)	30 (7%)	4 (1%)	21	55
All	All	1764/1844 (96%)	1633 (93%)	115 (6%)	16 (1%)	21	55

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	389	ASP
1	B	389	ASP
1	C	389	ASP
1	D	389	ASP
1	A	81	LYS
1	B	81	LYS
1	C	81	LYS
1	D	81	LYS
1	B	446	GLU
1	D	446	GLU
1	A	446	GLU
1	C	243	PRO
1	C	328	ASN
1	B	243	PRO
1	A	243	PRO
1	D	243	PRO

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	385/397 (97%)	358 (93%)	27 (7%)	19	47
1	B	385/397 (97%)	357 (93%)	28 (7%)	17	44
1	C	385/397 (97%)	359 (93%)	26 (7%)	20	49
1	D	385/397 (97%)	357 (93%)	28 (7%)	17	44
All	All	1540/1588 (97%)	1431 (93%)	109 (7%)	18	46

All (109) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	GLU
1	A	48	TRP
1	A	58	THR
1	A	78	ASN
1	A	84	LEU
1	A	97	LYS
1	A	199	GLU
1	A	220	GLN
1	A	234	ARG
1	A	241	LYS
1	A	244	GLU
1	A	247	ASP
1	A	255	VAL
1	A	267	SER
1	A	269	LYS
1	A	277	THR
1	A	286	LEU
1	A	302	ASN
1	A	307	VAL
1	A	328	ASN
1	A	338	ARG
1	A	340	VAL
1	A	347	LEU
1	A	367	LEU
1	A	388	THR
1	A	400	LYS
1	A	403	LEU
1	B	38	GLU
1	B	48	TRP
1	B	58	THR

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Mol	Chain	Res	Type
1	B	78	ASN
1	B	84	LEU
1	B	97	LYS
1	B	199	GLU
1	B	220	GLN
1	B	234	ARG
1	B	241	LYS
1	B	244	GLU
1	B	247	ASP
1	B	255	VAL
1	B	267	SER
1	B	269	LYS
1	B	277	THR
1	B	286	LEU
1	B	302	ASN
1	B	307	VAL
1	B	328	ASN
1	B	338	ARG
1	B	340	VAL
1	B	347	LEU
1	B	367	LEU
1	B	388	THR
1	B	400	LYS
1	B	403	LEU
1	B	441	LEU
1	C	38	GLU
1	C	48	TRP
1	C	58	THR
1	C	78	ASN
1	C	84	LEU
1	C	97	LYS
1	C	199	GLU
1	C	220	GLN
1	C	234	ARG
1	C	241	LYS
1	C	244	GLU
1	C	247	ASP
1	C	249	ARG
1	C	255	VAL
1	C	267	SER
1	C	269	LYS
1	C	286	LEU

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Mol	Chain	Res	Type
1	C	307	VAL
1	C	328	ASN
1	C	338	ARG
1	C	347	LEU
1	C	367	LEU
1	C	388	THR
1	C	400	LYS
1	C	403	LEU
1	C	441	LEU
1	D	38	GLU
1	D	48	TRP
1	D	58	THR
1	D	78	ASN
1	D	84	LEU
1	D	97	LYS
1	D	199	GLU
1	D	220	GLN
1	D	234	ARG
1	D	241	LYS
1	D	244	GLU
1	D	247	ASP
1	D	255	VAL
1	D	267	SER
1	D	269	LYS
1	D	277	THR
1	D	286	LEU
1	D	302	ASN
1	D	307	VAL
1	D	328	ASN
1	D	338	ARG
1	D	340	VAL
1	D	347	LEU
1	D	367	LEU
1	D	388	THR
1	D	400	LYS
1	D	403	LEU
1	D	441	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	ASN

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Mol	Chain	Res	Type
1	A	78	ASN
1	A	128	ASN
1	A	148	GLN
1	A	220	GLN
1	A	328	ASN
1	A	419	HIS
1	B	63	ASN
1	B	78	ASN
1	B	128	ASN
1	B	133	HIS
1	B	148	GLN
1	B	220	GLN
1	B	328	ASN
1	C	63	ASN
1	C	78	ASN
1	C	128	ASN
1	C	148	GLN
1	C	220	GLN
1	C	328	ASN
1	D	63	ASN
1	D	128	ASN
1	D	148	GLN
1	D	220	GLN
1	D	328	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 24 ligands modelled in this entry, 4 are monoatomic - leaving 20 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	$\# Z > 2$
5	EPE	A	1452	-	14,15,15	0.42	0	18,20,20	2.56	7 (38%)
6	OXY	A	1453	-	1,1,1	0.42	0	0,0,0	0.00	-
6	OXY	A	1454	-	1,1,1	0.39	0	0,0,0	0.00	-
2	FAD	A	500	-	48,58,58	1.20	5 (10%)	54,89,89	2.11	8 (14%)
3	NAP	A	501	-	42,52,52	1.42	4 (9%)	54,80,80	2.13	10 (18%)
5	EPE	B	1452	-	14,15,15	0.50	0	18,20,20	2.28	7 (38%)
6	OXY	B	1453	-	1,1,1	0.39	0	0,0,0	0.00	-
6	OXY	B	1454	-	1,1,1	0.44	0	0,0,0	0.00	-
2	FAD	B	500	-	48,58,58	1.26	5 (10%)	54,89,89	1.98	11 (20%)
3	NAP	B	501	-	42,52,52	1.38	6 (14%)	54,80,80	2.05	11 (20%)
5	EPE	C	1451	-	14,15,15	0.46	0	18,20,20	2.34	6 (33%)
6	OXY	C	1452	-	1,1,1	0.40	0	0,0,0	0.00	-
6	OXY	C	1453	-	1,1,1	0.42	0	0,0,0	0.00	-
2	FAD	C	500	-	48,58,58	1.18	4 (8%)	54,89,89	2.03	10 (18%)
3	NAP	C	501	-	42,52,52	1.44	5 (11%)	54,80,80	1.87	11 (20%)
5	EPE	D	1452	-	14,15,15	0.33	0	18,20,20	2.22	5 (27%)
6	OXY	D	1453	-	1,1,1	0.41	0	0,0,0	0.00	-
6	OXY	D	1454	-	1,1,1	0.39	0	0,0,0	0.00	-
2	FAD	D	500	-	48,58,58	1.28	6 (12%)	54,89,89	2.05	11 (20%)
3	NAP	D	501	-	42,52,52	1.31	5 (11%)	54,80,80	1.91	10 (18%)

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
5	EPE	A	1452	-	-	0/9/19/19	0/1/1/1
6	OXY	A	1453	-	-	0/0/0/0	0/0/0/0
6	OXY	A	1454	-	-	0/0/0/0	0/0/0/0
2	FAD	A	500	-	-	0/30/50/50	0/6/6/6
3	NAP	A	501	-	-	0/27/67/67	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EPE	B	1452	-	-	0/9/19/19	0/1/1/1
6	OXY	B	1453	-	-	0/0/0/0	0/0/0/0
6	OXY	B	1454	-	-	0/0/0/0	0/0/0/0
2	FAD	B	500	-	-	0/30/50/50	0/6/6/6
3	NAP	B	501	-	-	0/27/67/67	0/5/5/5
5	EPE	C	1451	-	-	0/9/19/19	0/1/1/1
6	OXY	C	1452	-	-	0/0/0/0	0/0/0/0
6	OXY	C	1453	-	-	0/0/0/0	0/0/0/0
2	FAD	C	500	-	-	0/30/50/50	0/6/6/6
3	NAP	C	501	-	-	0/27/67/67	0/5/5/5
5	EPE	D	1452	-	-	0/9/19/19	0/1/1/1
6	OXY	D	1453	-	-	0/0/0/0	0/0/0/0
6	OXY	D	1454	-	-	0/0/0/0	0/0/0/0
2	FAD	D	500	-	-	0/30/50/50	0/6/6/6
3	NAP	D	501	-	-	0/27/67/67	0/5/5/5

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	NAP	C7N-N7N	2.02	1.37	1.33
3	B	501	NAP	PA-O2A	2.02	1.63	1.54
3	B	501	NAP	P2B-O3X	2.03	1.62	1.54
2	D	500	FAD	C10-N1	2.05	1.39	1.35
2	A	500	FAD	C2A-N1A	2.05	1.37	1.33
2	C	500	FAD	C1'-N10	2.08	1.50	1.48
3	D	501	NAP	P2B-O2X	2.10	1.62	1.54
3	A	501	NAP	C7N-N7N	2.10	1.37	1.33
3	D	501	NAP	C4N-C3N	2.16	1.43	1.39
3	C	501	NAP	C7N-N7N	2.20	1.37	1.33
2	A	500	FAD	C1'-N10	2.22	1.50	1.48
2	A	500	FAD	C4-N3	2.23	1.37	1.33
2	D	500	FAD	C1'-N10	2.24	1.50	1.48
3	C	501	NAP	P2B-O3X	2.30	1.63	1.54
3	B	501	NAP	P2B-O2X	2.31	1.63	1.54
3	C	501	NAP	P2B-O2X	2.32	1.63	1.54
3	A	501	NAP	P2B-O2X	2.37	1.63	1.54
3	D	501	NAP	C2N-C3N	2.38	1.42	1.39
2	B	500	FAD	C1'-N10	2.39	1.50	1.48
2	B	500	FAD	C2A-N1A	2.40	1.38	1.33
2	D	500	FAD	C2A-N1A	2.40	1.38	1.33
2	D	500	FAD	C4-N3	2.51	1.37	1.33
2	A	500	FAD	C4X-N5	2.59	1.37	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	500	FAD	C4-N3	3.12	1.38	1.33
2	B	500	FAD	C4-N3	3.15	1.39	1.33
2	B	500	FAD	C4X-N5	3.22	1.38	1.33
2	D	500	FAD	C4X-N5	3.26	1.38	1.33
2	C	500	FAD	C4X-N5	3.28	1.38	1.33
2	B	500	FAD	C2A-N3A	3.29	1.38	1.32
2	D	500	FAD	C2A-N3A	3.39	1.38	1.32
2	A	500	FAD	C2A-N3A	3.41	1.38	1.32
2	C	500	FAD	C2A-N3A	3.42	1.38	1.32
3	D	501	NAP	O4B-C1B	3.56	1.45	1.41
3	D	501	NAP	O4D-C1D	3.61	1.45	1.41
3	B	501	NAP	O4B-C1B	3.64	1.45	1.41
3	C	501	NAP	O4B-C1B	3.66	1.45	1.41
3	A	501	NAP	O4D-C1D	3.70	1.45	1.41
3	A	501	NAP	O4B-C1B	4.06	1.46	1.41
3	C	501	NAP	O4D-C1D	4.95	1.47	1.41
3	B	501	NAP	O4D-C1D	5.03	1.47	1.41

All (107) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	N3A-C2A-N1A	-10.19	121.09	128.89
2	C	500	FAD	N3A-C2A-N1A	-9.82	121.37	128.89
2	D	500	FAD	N3A-C2A-N1A	-9.67	121.49	128.89
2	B	500	FAD	N3A-C2A-N1A	-8.40	122.47	128.89
3	A	501	NAP	N3A-C2A-N1A	-8.14	122.66	128.89
3	B	501	NAP	N3A-C2A-N1A	-7.53	123.13	128.89
3	D	501	NAP	N3A-C2A-N1A	-7.06	123.49	128.89
3	B	501	NAP	P2B-O2B-C2B	-5.47	108.44	121.56
3	D	501	NAP	PN-O3-PA	-5.46	117.39	132.73
3	A	501	NAP	PN-O3-PA	-5.22	118.07	132.73
3	B	501	NAP	PN-O3-PA	-5.19	118.16	132.73
3	C	501	NAP	PN-O3-PA	-5.18	118.18	132.73
3	A	501	NAP	P2B-O2B-C2B	-4.89	109.84	121.56
3	C	501	NAP	N3A-C2A-N1A	-4.84	125.19	128.89
3	C	501	NAP	P2B-O2B-C2B	-4.51	110.76	121.56
2	A	500	FAD	P-O3P-PA	-4.39	120.41	132.73
3	D	501	NAP	P2B-O2B-C2B	-4.04	111.88	121.56
2	B	500	FAD	P-O3P-PA	-3.81	122.03	132.73
2	C	500	FAD	P-O3P-PA	-3.59	122.64	132.73
2	D	500	FAD	P-O3P-PA	-3.45	123.03	132.73
3	A	501	NAP	C4B-O4B-C1B	-3.44	105.94	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500	FAD	C4A-C5A-N7A	-3.28	106.47	109.48
2	B	500	FAD	C4X-C4-N3	-3.04	119.44	123.59
2	A	500	FAD	C4A-C5A-N7A	-2.89	106.82	109.48
3	C	501	NAP	C3B-C2B-C1B	-2.76	97.39	102.73
3	A	501	NAP	C4A-C5A-N7A	-2.73	106.96	109.48
3	B	501	NAP	C4A-C5A-N7A	-2.71	106.99	109.48
2	B	500	FAD	C4A-C5A-N7A	-2.70	106.99	109.48
3	C	501	NAP	C5N-C4N-C3N	-2.62	117.04	120.33
3	C	501	NAP	C4B-O4B-C1B	-2.60	106.86	109.72
2	D	500	FAD	C4A-C5A-N7A	-2.57	107.12	109.48
3	A	501	NAP	C5N-C4N-C3N	-2.53	117.15	120.33
2	A	500	FAD	O2'-C2'-C1'	-2.48	103.86	109.94
2	B	500	FAD	O2'-C2'-C1'	-2.45	103.91	109.94
3	D	501	NAP	O7N-C7N-N7N	-2.40	119.22	122.59
3	B	501	NAP	O7N-C7N-N7N	-2.33	119.31	122.59
3	D	501	NAP	C4B-O4B-C1B	-2.32	107.17	109.72
2	D	500	FAD	O2'-C2'-C1'	-2.25	104.41	109.94
2	C	500	FAD	C4X-C4-N3	-2.24	120.53	123.59
2	C	500	FAD	C1B-N9A-C4A	-2.14	123.71	126.94
2	D	500	FAD	C9A-C5X-N5	-2.10	119.25	122.36
3	D	501	NAP	C4A-C5A-N7A	-2.08	107.56	109.48
3	C	501	NAP	C4A-C5A-N7A	-2.05	107.60	109.48
3	B	501	NAP	C3B-C2B-C1B	-2.04	98.77	102.73
2	D	500	FAD	C4X-C4-N3	-2.04	120.80	123.59
2	C	500	FAD	O2P-P-O3P	2.00	114.17	105.09
2	B	500	FAD	C5X-C9A-N10	2.00	119.14	117.62
3	B	501	NAP	O3-PA-O5B	2.06	108.41	102.94
5	B	1452	EPE	C2-C3-N4	2.11	114.41	110.63
2	D	500	FAD	C6-C5X-C9A	2.16	121.83	118.98
2	A	500	FAD	C6-C5X-C9A	2.17	121.84	118.98
3	B	501	NAP	O4B-C1B-N9A	2.23	112.77	108.10
5	A	1452	EPE	C6-N1-C2	2.28	113.84	108.90
3	A	501	NAP	O4D-C1D-N1N	2.29	110.65	108.13
5	C	1451	EPE	C2-C3-N4	2.32	114.78	110.63
5	B	1452	EPE	C6-N1-C2	2.32	113.93	108.90
2	B	500	FAD	O3P-PA-O5B	2.34	109.14	102.94
5	A	1452	EPE	C7-N4-C5	2.38	117.36	111.27
5	B	1452	EPE	C7-N4-C5	2.39	117.39	111.27
5	D	1452	EPE	C2-C3-N4	2.44	115.00	110.63
2	B	500	FAD	C6-C5X-C9A	2.50	122.27	118.98
5	C	1451	EPE	C6-N1-C2	2.54	114.41	108.90
2	D	500	FAD	C5X-C9A-N10	2.57	119.58	117.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	1452	EPE	O2S-S-C10	2.60	109.12	106.91
5	D	1452	EPE	C7-N4-C5	2.60	117.92	111.27
5	D	1452	EPE	C7-N4-C3	2.63	118.00	111.27
5	C	1451	EPE	C6-C5-N4	2.65	115.37	110.63
5	A	1452	EPE	C6-C5-N4	2.67	115.41	110.63
2	C	500	FAD	C4X-N5-C5X	2.72	119.89	116.76
5	A	1452	EPE	C7-N4-C3	2.74	118.30	111.27
3	D	501	NAP	O4B-C1B-N9A	2.74	113.84	108.10
3	A	501	NAP	O4B-C1B-N9A	2.78	113.91	108.10
3	D	501	NAP	C3N-C7N-N7N	2.79	120.87	117.82
2	C	500	FAD	C5X-C9A-N10	2.81	119.76	117.62
3	C	501	NAP	O4B-C1B-N9A	2.84	114.05	108.10
5	B	1452	EPE	C7-N4-C3	2.90	118.71	111.27
5	C	1451	EPE	C7-N4-C3	3.01	118.99	111.27
5	B	1452	EPE	C6-C5-N4	3.06	116.10	110.63
3	D	501	NAP	O4D-C1D-N1N	3.22	111.67	108.13
3	B	501	NAP	C3N-C7N-N7N	3.25	121.38	117.82
5	B	1452	EPE	O2S-S-C10	3.29	109.71	106.91
2	B	500	FAD	C1'-N10-C9A	3.38	122.65	118.86
5	A	1452	EPE	O2S-S-C10	3.46	109.86	106.91
2	A	500	FAD	C4X-N5-C5X	3.64	120.95	116.76
3	C	501	NAP	C3N-C7N-N7N	3.70	121.87	117.82
3	C	501	NAP	O3-PN-O5D	3.71	112.79	102.94
2	D	500	FAD	C4X-N5-C5X	3.91	121.25	116.76
3	B	501	NAP	O4D-C1D-N1N	3.92	112.43	108.13
3	D	501	NAP	O3-PN-O5D	3.93	113.37	102.94
2	C	500	FAD	C1'-N10-C9A	4.06	123.42	118.86
2	B	500	FAD	C4X-N5-C5X	4.08	121.45	116.76
3	A	501	NAP	C3N-C7N-N7N	4.13	122.34	117.82
2	A	500	FAD	C1'-N10-C9A	4.32	123.72	118.86
3	B	501	NAP	O3-PN-O5D	4.38	114.55	102.94
2	D	500	FAD	C1'-N10-C9A	4.44	123.85	118.86
3	C	501	NAP	O4D-C1D-N1N	4.55	113.13	108.13
2	A	500	FAD	C4-N3-C2	4.92	119.50	115.25
2	C	500	FAD	C4-N3-C2	4.97	119.54	115.25
2	D	500	FAD	C4-N3-C2	5.00	119.57	115.25
5	C	1451	EPE	O2S-S-C10	5.04	111.21	106.91
5	C	1451	EPE	C5-N4-C3	5.19	120.14	108.90
3	A	501	NAP	O3-PN-O5D	5.24	116.84	102.94
2	B	500	FAD	C4-N3-C2	5.81	120.27	115.25
5	B	1452	EPE	C5-N4-C3	5.91	121.70	108.90
5	A	1452	EPE	O1S-S-C10	5.98	112.01	106.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1452	EPE	C5-N4-C3	6.04	121.97	108.90
5	D	1452	EPE	C5-N4-C3	6.87	123.78	108.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1452	EPE	4	0
2	A	500	FAD	1	0
3	A	501	NAP	2	0
5	B	1452	EPE	5	0
2	B	500	FAD	1	0
3	B	501	NAP	3	0
3	C	501	NAP	2	0
2	D	500	FAD	1	0
3	D	501	NAP	2	0

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 ⓘ

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	443/461 (96%)	-0.22	5 (1%) 82 74	40, 52, 67, 74	0
1	B	443/461 (96%)	-0.15	6 (1%) 78 69	40, 52, 67, 74	0
1	C	443/461 (96%)	-0.20	5 (1%) 82 74	40, 52, 67, 74	0
1	D	443/461 (96%)	-0.25	3 (0%) 89 84	40, 52, 67, 74	0
All	All	1772/1844 (96%)	-0.21	19 (1%) 82 74	40, 52, 67, 74	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	244	GLU	4.2
1	A	236	ALA	3.1
1	A	235	THR	2.9
1	B	139	ASP	2.7
1	B	241	LYS	2.6
1	B	450	SER	2.5
1	D	244	GLU	2.4
1	C	446	GLU	2.4
1	D	139	ASP	2.4
1	C	328	ASN	2.4
1	A	368	VAL	2.3
1	D	351	GLU	2.3
1	B	446	GLU	2.2
1	A	446	GLU	2.2
1	C	256	ASP	2.2
1	C	326	SER	2.1
1	A	328	ASN	2.1
1	B	244	GLU	2.0
1	B	240	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, 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	LLDF	B-factors(Å ²)	Q<0.9
6	OXY	D	1454	2/2	0.95	0.69	9.50	52,52,52,54	0
5	EPE	A	1452	15/15	0.86	0.24	4.68	97,103,107,108	0
5	EPE	C	1451	15/15	0.77	0.30	4.55	102,104,110,111	0
5	EPE	D	1452	15/15	0.82	0.26	4.46	103,104,107,107	0
5	EPE	B	1452	15/15	0.84	0.23	2.51	106,107,108,108	0
6	OXY	B	1453	2/2	0.91	0.19	1.37	68,68,68,68	0
6	OXY	A	1454	2/2	0.93	0.19	1.22	55,55,55,55	0
3	NAP	C	501	48/48	0.93	0.20	0.48	41,44,50,52	0
3	NAP	A	501	48/48	0.94	0.19	0.26	41,44,50,51	0
2	FAD	D	500	53/53	0.97	0.20	0.25	37,41,46,48	0
3	NAP	B	501	48/48	0.94	0.18	0.15	41,44,50,51	0
2	FAD	A	500	53/53	0.97	0.19	-0.07	37,41,46,48	0
2	FAD	C	500	53/53	0.97	0.18	-0.24	37,40,46,48	0
2	FAD	B	500	53/53	0.97	0.18	-0.28	37,41,46,48	0
3	NAP	D	501	48/48	0.94	0.16	-0.38	41,44,50,51	0
6	OXY	B	1454	2/2	0.94	0.65	-	53,53,53,54	0
6	OXY	D	1453	2/2	0.96	0.15	-	45,45,45,46	0
4	CL	D	1451	1/1	0.97	0.09	-	56,56,56,56	0
6	OXY	C	1452	2/2	0.90	0.25	-	50,50,50,51	0
4	CL	B	1451	1/1	0.91	0.12	-	60,60,60,60	0
4	CL	A	1451	1/1	0.98	0.12	-	52,52,52,52	0
4	CL	C	1454	1/1	0.98	0.15	-	58,58,58,58	0
6	OXY	A	1453	2/2	0.97	0.58	-	49,49,49,50	0
6	OXY	C	1453	2/2	0.96	0.58	-	51,51,51,51	0

6.5 Other polymers

There are no such residues in this entry.