



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 02:02 PM UTC

PDB ID : 5VPN / pdb_00005vpn
Title : E. coli Quinol fumarate reductase FrdA E245Q mutation
Authors : Starbird, C.A.; Maklashina, E.; Sharma, P.; Qualls-Histed, S.; Cecchini, G.; Iverson, T.M.
Deposited on : 2017-05-05
Resolution : 4.22 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

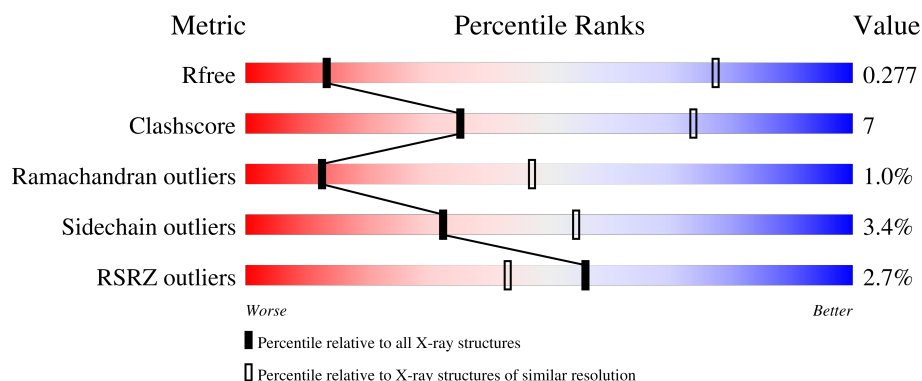
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.22 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1014 (4.56-3.88)
Clashscore	190562	1033 (4.54-3.90)
Ramachandran outliers	187476	1002 (4.58-3.86)
Sidechain outliers	187428	1030 (4.60-3.84)
RSRZ outliers	180081	1011 (4.56-3.88)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	585	
1	E	585	
2	B	243	
2	F	243	
3	C	130	

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Mol	Chain	Length	Quality of chain
3	G	130	% 77% 21%
4	D	119	79% 18%
4	H	119	82% 15%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 16845 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 Fumarate reductase flavoprotein subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	577	Total	C	N	O	S	0	0	0
			4448	2775	803	839	31			
1	E	585	Total	C	N	O	S	0	0	0
			4509	2813	815	850	31			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	245	GLN	GLU	engineered mutation	UNP P00363
E	245	GLN	GLU	engineered mutation	UNP P00363

- Molecule 2 is a protein called Succinate dehydrogenase iron-sulfur subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	243	Total	C	N	O	S	0	0	0
			1888	1189	323	357	19			
2	F	243	Total	C	N	O	S	0	0	0
			1888	1189	323	357	19			

- Molecule 3 is a protein called Fumarate reductase subunit C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	130	Total	C	N	O	S	0	0	0
			1058	720	166	169	3			
3	G	130	Total	C	N	O	S	0	0	0
			1058	720	166	169	3			

- Molecule 4 is a protein called Fumarate reductase subunit D.

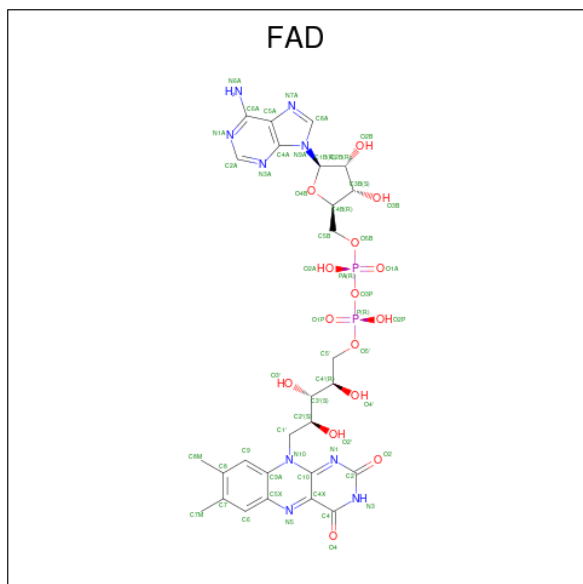
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	119	Total	C	N	O	S	0	0	0
			926	626	151	142	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	119	Total	C	N	O	S	0	0	0
			926	626	151	142	7			

- Molecule 5 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



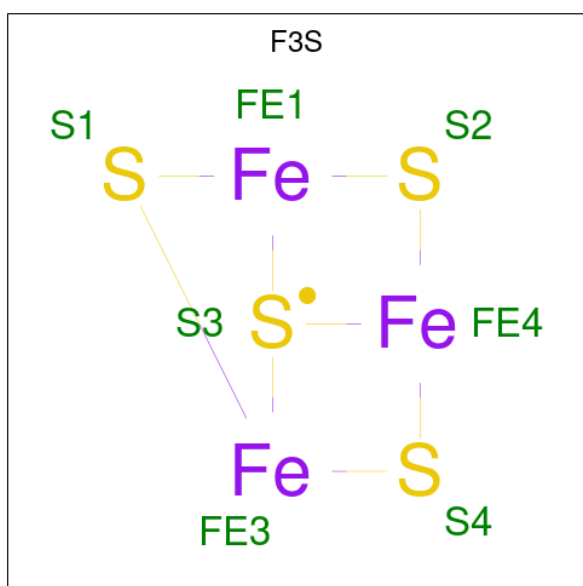
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
5	E	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 6 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



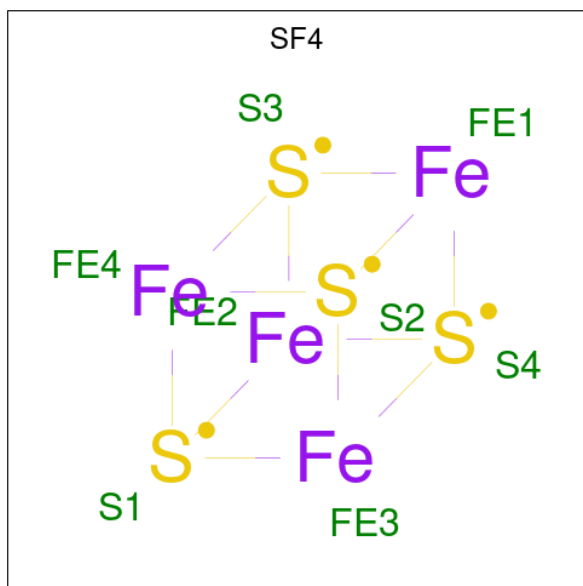
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	Fe	S	0	0
			4	2	2		
6	F	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 7 is FE3-S4 CLUSTER (CCD ID: F3S) (formula: Fe₃S₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	Fe	S	0	0
			7	3	4		
7	F	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 8 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).

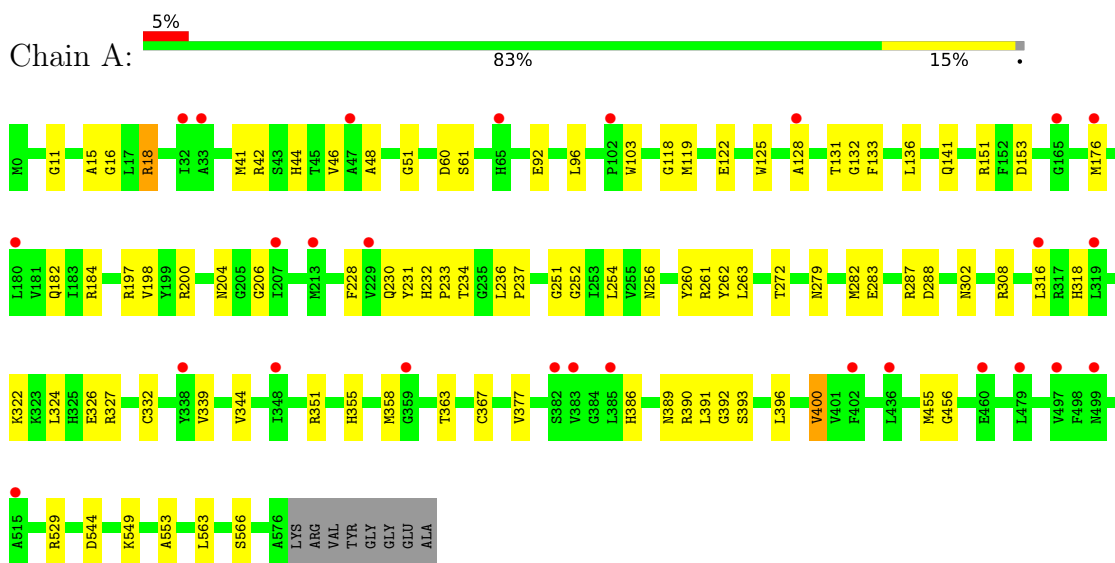


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	Fe	S	0	0
			8	4	4		
8	F	1	Total	Fe	S	0	0
			8	4	4		

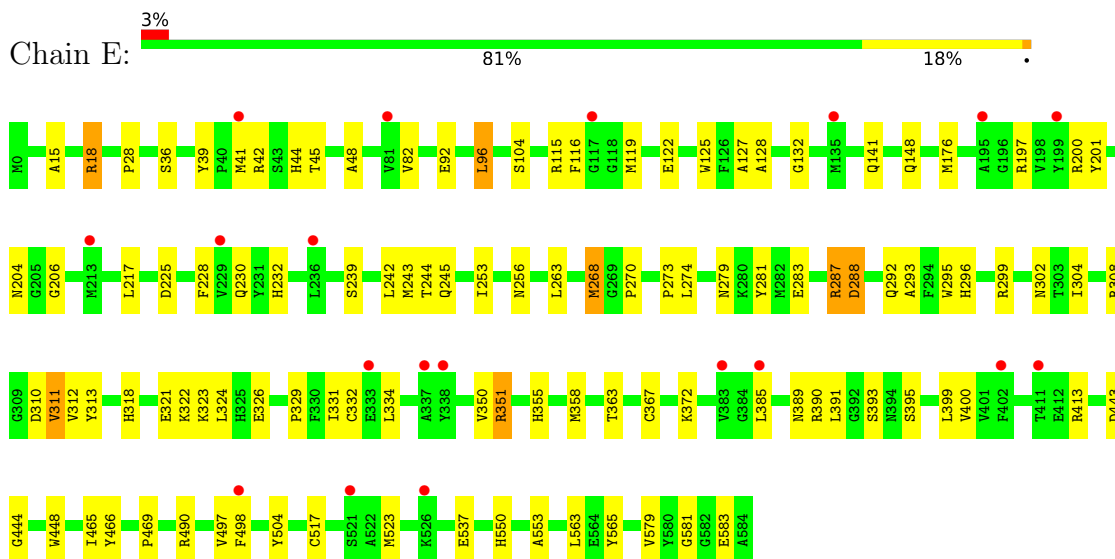
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

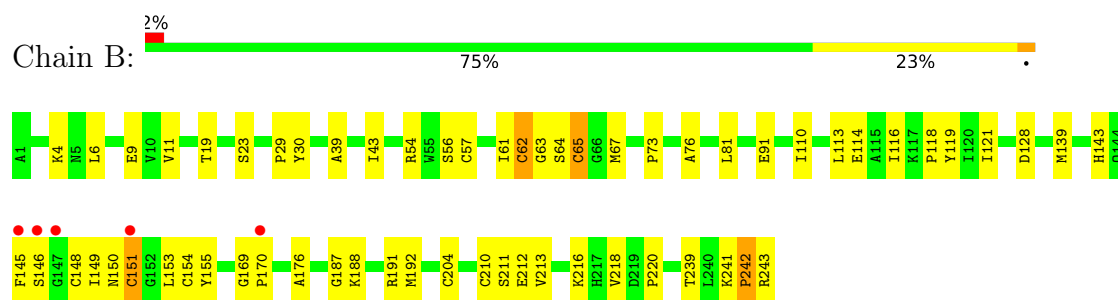
- Molecule 1: Fumarate reductase flavoprotein subunit



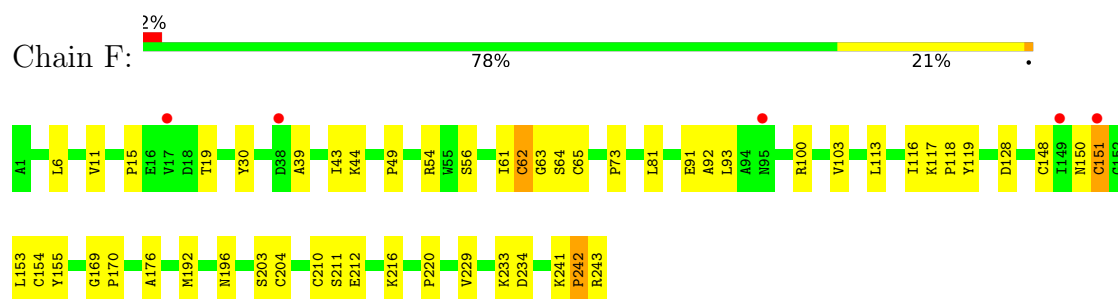
- Molecule 1: Fumarate reductase flavoprotein subunit



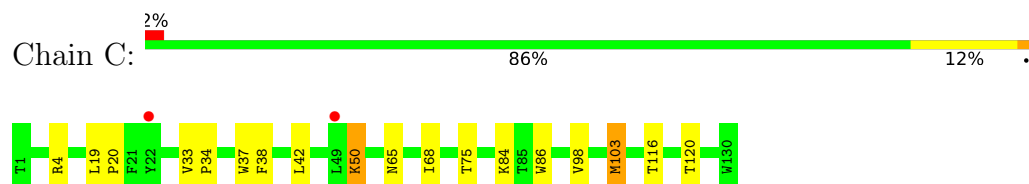
- Molecule 2: Succinate dehydrogenase iron-sulfur subunit



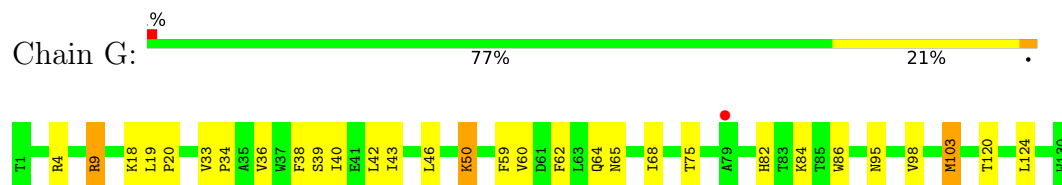
• Molecule 2: Succinate dehydrogenase iron-sulfur subunit



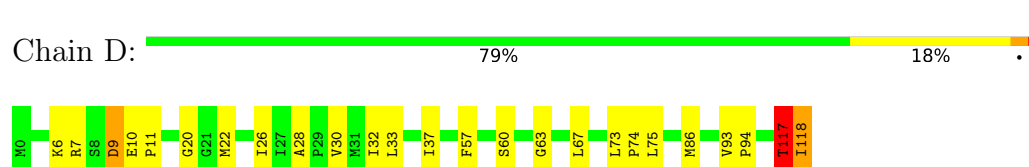
• Molecule 3: Fumarate reductase subunit C



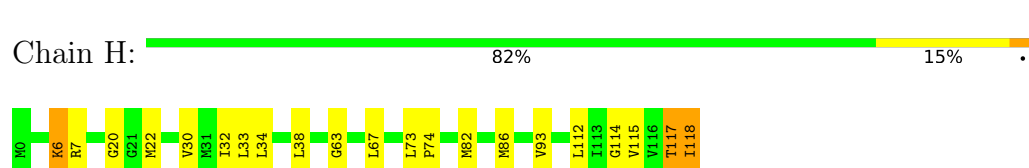
• Molecule 3: Fumarate reductase subunit C



• Molecule 4: Fumarate reductase subunit D



• Molecule 4: Fumarate reductase subunit D



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	133.56Å 138.13Å 220.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.96 – 4.22 22.96 – 4.22	Depositor EDS
% Data completeness (in resolution range)	95.9 (22.96-4.22) 80.7 (22.96-4.22)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.24 (at 4.14Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.215 , 0.270 0.221 , 0.277	Depositor DCC
R_{free} test set	1437 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	135.1	Xtriage
Anisotropy	0.680	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 146.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.017 for k,h,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	16845	wwPDB-VP
Average B, all atoms (Å ²)	214.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FES, SF4, FAD, F3S

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	0.16	0/4540	0.36	0/6139
1	E	0.17	0/4602	0.39	0/6221
2	B	0.18	0/1931	0.40	0/2617
2	F	0.16	0/1931	0.38	0/2617
3	C	0.17	0/1094	0.39	0/1496
3	G	0.18	0/1094	0.43	0/1496
4	D	0.19	0/956	0.38	0/1303
4	H	0.21	0/956	0.43	0/1303
All	All	0.17	0/17104	0.38	0/23192

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	4448	0	4337	60	0
1	E	4509	0	4398	79	0
2	B	1888	0	1848	38	0
2	F	1888	0	1847	36	0
3	C	1058	0	1108	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	1058	0	1108	25	0
4	D	926	0	971	19	0
4	H	926	0	971	16	0
5	A	53	0	31	3	0
5	E	53	0	31	10	0
6	B	4	0	0	0	0
6	F	4	0	0	0	0
7	B	7	0	0	0	0
7	F	7	0	0	0	0
8	B	8	0	0	0	0
8	F	8	0	0	0	0
All	All	16845	0	16650	249	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 (249) 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:E:96:LEU:HD11	1:E:400:VAL:HG11	1.54	0.87
3:G:65:ASN:HB3	3:G:68:ILE:H	1.50	0.76
2:B:62:CYS:SG	2:B:63:GLY:N	2.60	0.74
1:E:28:PRO:HA	1:E:148:GLN:HE21	1.52	0.74
1:E:41:MET:HE2	2:F:150:ASN:HD22	1.51	0.74
1:E:176:MET:HB3	3:G:4:ARG:HD3	1.69	0.74
1:A:96:LEU:HD11	1:A:400:VAL:HG11	1.71	0.72
3:C:65:ASN:HB3	3:C:68:ILE:H	1.55	0.72
4:D:20:GLY:HA2	4:D:73:LEU:HB3	1.72	0.71
1:A:288:ASP:OD1	1:A:389:ASN:ND2	2.23	0.71
1:A:324:LEU:HD13	1:A:344:VAL:HA	1.71	0.70
3:C:98:VAL:HG22	3:C:103:MET:HB3	1.74	0.70
3:G:19:LEU:HD12	3:G:20:PRO:HD2	1.73	0.70
1:E:293:ALA:HA	1:E:296:HIS:HD1	1.58	0.69
1:E:393:SER:HA	5:E:601:FAD:HN3	1.59	0.68
1:E:197:ARG:HD2	1:E:206:GLY:HA2	1.77	0.67
1:E:48:ALA:HB3	1:E:132:GLY:HA3	1.77	0.67
1:E:281:TYR:HE1	1:E:583:GLU:HB2	1.59	0.66
1:E:395:SER:HG	5:E:601:FAD:HO3'	1.44	0.66
1:A:197:ARG:HD2	1:A:206:GLY:HA2	1.79	0.64
2:F:93:LEU:O	3:G:9:ARG:NH2	2.31	0.64
1:A:151:ARG:NH1	1:A:153:ASP:OD2	2.31	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:120:THR:HG23	4:D:30:VAL:HB	1.79	0.63
1:E:270:PRO:HD2	1:E:281:TYR:HE2	1.65	0.62
2:F:44:LYS:NZ	2:F:49:PRO:O	2.31	0.62
1:E:18:ARG:NH1	1:E:92:GLU:OE1	2.34	0.61
1:A:529:ARG:NH2	1:A:544:ASP:OD1	2.32	0.60
4:H:34:LEU:HD23	4:H:38:LEU:HD12	1.82	0.60
3:G:75:THR:HG22	4:H:32:ILE:HD13	1.82	0.59
1:E:45:THR:OG1	5:E:601:FAD:O2A	2.19	0.59
4:H:20:GLY:HA2	4:H:73:LEU:HB3	1.84	0.59
3:C:33:VAL:HB	3:C:34:PRO:HD3	1.84	0.59
1:A:251:GLY:O	1:A:318:HIS:NE2	2.29	0.59
3:C:50:LYS:HB3	4:D:118:ILE:HG22	1.84	0.58
3:G:120:THR:HG23	4:H:30:VAL:HB	1.85	0.58
1:A:18:ARG:NH1	1:A:92:GLU:OE1	2.36	0.58
1:E:281:TYR:CE1	1:E:583:GLU:HB2	2.38	0.58
2:F:62:CYS:SG	2:F:63:GLY:N	2.77	0.58
1:A:230:GLN:OE1	1:A:390:ARG:NH2	2.35	0.58
1:A:18:ARG:HG2	1:A:400:VAL:HA	1.86	0.58
2:F:54:ARG:HB3	2:F:64:SER:HB2	1.87	0.57
3:G:98:VAL:HG22	3:G:103:MET:HB3	1.87	0.57
1:A:11:GLY:HA2	5:A:601:FAD:H1B	1.87	0.57
2:F:192:MET:O	2:F:196:ASN:ND2	2.38	0.57
1:E:390:ARG:HD2	1:E:395:SER:HB2	1.86	0.57
1:E:256:ASN:HD22	1:E:302:ASN:HB3	1.68	0.56
1:A:390:ARG:NH1	1:A:391:LEU:O	2.38	0.56
2:B:54:ARG:HB3	2:B:64:SER:HB2	1.86	0.56
3:C:75:THR:HG22	4:D:32:ILE:HD13	1.86	0.56
1:E:355:HIS:HE1	5:E:601:FAD:C8	2.19	0.56
2:F:11:VAL:HG21	2:F:91:GLU:HG2	1.88	0.56
2:B:143:HIS:O	2:B:146:SER:OG	2.23	0.55
2:F:6:LEU:HD23	2:F:81:LEU:HD13	1.88	0.55
2:B:30:TYR:HB3	2:B:81:LEU:HD12	1.89	0.55
2:F:30:TYR:HB3	2:F:81:LEU:HD12	1.88	0.55
2:F:211:SER:HA	2:F:220:PRO:HD2	1.89	0.54
1:E:292:GLN:HA	1:E:465:ILE:HG21	1.89	0.54
3:G:33:VAL:HB	3:G:34:PRO:HD3	1.90	0.54
1:A:237:PRO:HB2	1:A:308:ARG:HB2	1.89	0.54
2:B:210:CYS:SG	2:B:220:PRO:HG2	2.48	0.54
1:E:288:ASP:CG	1:E:389:ASN:HD21	2.16	0.53
1:E:393:SER:HA	5:E:601:FAD:N3	2.23	0.53
1:A:393:SER:HA	5:A:601:FAD:HN3	1.72	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:73:PRO:HB2	2:F:153:LEU:HD22	1.90	0.53
1:E:295:TRP:HE1	1:E:299:ARG:NE	2.07	0.53
1:A:176:MET:HG3	3:C:4:ARG:HD3	1.91	0.53
1:E:44:HIS:CE1	1:E:204:ASN:HA	2.43	0.53
1:E:82:VAL:HG22	1:E:385:LEU:HD12	1.91	0.53
2:F:155:TYR:CE2	2:F:169:GLY:HA3	2.43	0.53
1:E:253:ILE:HG12	1:E:318:HIS:HE1	1.74	0.53
1:E:141:GLN:HB3	2:F:118:PRO:O	2.08	0.53
1:E:232:HIS:HB2	1:E:355:HIS:HD2	1.73	0.53
2:F:234:ASP:OD1	4:H:7:ARG:NH2	2.42	0.52
2:B:11:VAL:HG21	2:B:91:GLU:HG2	1.92	0.52
4:D:6:LYS:HE2	4:H:6:LYS:HB2	1.90	0.52
2:F:210:CYS:SG	2:F:220:PRO:HG2	2.50	0.52
3:G:86:TRP:HE1	4:H:22:MET:HE2	1.74	0.52
1:E:263:LEU:HD22	1:E:268:MET:HE3	1.91	0.52
3:G:46:LEU:HD21	4:H:114:GLY:HA3	1.92	0.52
3:C:50:LYS:HG3	4:D:118:ILE:H	1.75	0.51
1:A:254:LEU:HD12	1:A:263:LEU:HD11	1.91	0.51
2:F:151:CYS:SG	2:F:153:LEU:HG	2.50	0.51
1:A:393:SER:HA	5:A:601:FAD:N3	2.25	0.51
1:E:295:TRP:CD1	1:E:466:TYR:HD1	2.29	0.51
2:B:154:CYS:SG	2:B:170:PRO:HG2	2.51	0.51
1:E:232:HIS:HB2	1:E:355:HIS:CD2	2.46	0.51
1:A:41:MET:HE2	2:B:150:ASN:HD22	1.74	0.51
2:B:149:ILE:HG13	2:B:151:CYS:HB3	1.92	0.51
4:D:94:PRO:HA	2:F:242:PRO:HA	1.91	0.51
1:E:104:SER:OG	1:E:127:ALA:O	2.26	0.50
2:F:210:CYS:SG	2:F:211:SER:N	2.84	0.50
1:A:51:GLY:HA2	1:A:131:THR:HG21	1.93	0.50
2:B:151:CYS:SG	2:B:153:LEU:HG	2.51	0.50
3:C:19:LEU:HD12	3:C:20:PRO:HD2	1.93	0.50
1:E:18:ARG:HG2	1:E:400:VAL:HA	1.92	0.50
1:A:15:ALA:HB1	1:A:377:VAL:HG13	1.93	0.50
1:E:310:ASP:O	1:E:351:ARG:NH1	2.42	0.50
1:A:42:ARG:NH2	2:B:63:GLY:O	2.45	0.49
2:B:155:TYR:CZ	2:B:169:GLY:HA3	2.46	0.49
3:G:50:LYS:HB3	4:H:118:ILE:HG22	1.95	0.49
3:C:116:THR:HA	4:D:26:ILE:HG23	1.94	0.49
4:D:73:LEU:HB2	4:D:74:PRO:HD3	1.95	0.49
2:B:210:CYS:SG	2:B:211:SER:N	2.86	0.49
1:E:232:HIS:HE1	1:E:245:GLN:HG3	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:390:ARG:HH12	1:A:392:GLY:HA2	1.77	0.49
1:E:115:ARG:HD3	1:E:279:ASN:ND2	2.28	0.49
1:A:237:PRO:HG2	1:A:308:ARG:HD2	1.94	0.49
1:E:395:SER:N	5:E:601:FAD:O2	2.45	0.49
2:B:155:TYR:CE2	2:B:169:GLY:HA3	2.47	0.49
3:G:38:PHE:O	3:G:42:LEU:HG	2.13	0.48
4:D:33:LEU:HA	4:D:37:ILE:HD12	1.95	0.48
3:C:86:TRP:HE1	4:D:22:MET:HE2	1.79	0.48
1:E:116:PHE:CE2	1:E:245:GLN:HB2	2.49	0.48
1:E:372:LYS:HE3	1:E:413:ARG:HD2	1.95	0.47
1:E:497:VAL:HG21	2:F:15:PRO:HG2	1.96	0.47
1:E:331:ILE:HD13	1:E:334:LEU:HD12	1.96	0.47
1:E:119:MET:HE2	1:E:391:LEU:HD23	1.97	0.47
2:F:92:ALA:O	3:G:9:ARG:NH1	2.36	0.47
1:A:48:ALA:HB3	1:A:132:GLY:HA3	1.95	0.47
1:A:236:LEU:HD13	1:A:339:VAL:HG13	1.96	0.47
3:C:38:PHE:O	3:C:42:LEU:HG	2.14	0.47
3:G:36:VAL:O	3:G:40:ILE:HG12	2.15	0.47
1:E:253:ILE:HA	1:E:283:GLU:HG2	1.97	0.47
2:F:154:CYS:SG	2:F:170:PRO:HG2	2.55	0.47
3:G:50:LYS:HD2	4:H:118:ILE:O	2.15	0.46
2:B:145:PHE:HA	2:B:218:VAL:HG13	1.97	0.46
1:E:497:VAL:HG13	3:G:4:ARG:HG2	1.96	0.46
1:E:200:ARG:NH1	1:E:201:TYR:OH	2.48	0.46
2:B:6:LEU:HB2	2:B:30:TYR:CE2	2.51	0.46
4:D:57:PHE:O	4:D:60:SER:OG	2.32	0.46
3:G:59:PHE:O	3:G:62:PHE:HB3	2.15	0.46
1:A:322:LYS:O	1:A:326:GLU:HB3	2.15	0.46
3:G:60:VAL:O	3:G:64:GLN:HG3	2.15	0.46
1:A:141:GLN:HB3	2:B:118:PRO:O	2.16	0.46
2:B:73:PRO:HG2	2:B:213:VAL:HG11	1.96	0.46
1:E:295:TRP:CE3	1:E:465:ILE:HB	2.51	0.46
1:E:448:TRP:CH2	1:E:504:TYR:HB3	2.51	0.46
3:G:50:LYS:O	3:G:50:LYS:HD3	2.16	0.46
4:D:9:ASP:OD1	4:D:9:ASP:N	2.36	0.45
1:A:42:ARG:HG3	2:B:62:CYS:SG	2.56	0.45
2:B:67:MET:SD	2:B:76:ALA:HB2	2.56	0.45
1:A:252:GLY:HA3	1:A:316:LEU:HD23	1.96	0.45
2:B:113:LEU:HD23	2:B:118:PRO:HG3	1.97	0.45
2:B:188:LYS:O	2:B:192:MET:HG2	2.16	0.45
1:E:48:ALA:HA	5:E:601:FAD:C6	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:253:ILE:HG12	1:E:318:HIS:CE1	2.52	0.45
4:H:73:LEU:HB2	4:H:74:PRO:HD3	1.98	0.45
3:C:50:LYS:HD2	4:D:117:THR:HG23	1.98	0.45
1:E:230:GLN:CD	1:E:287:ARG:HD3	2.42	0.45
1:E:443:ASP:HA	1:E:490:ARG:HG3	1.97	0.45
1:A:11:GLY:O	1:A:16:GLY:HA3	2.16	0.45
1:E:15:ALA:HB2	1:E:399:LEU:HD22	1.99	0.45
1:E:550:HIS:O	1:E:565:TYR:HA	2.17	0.45
3:G:39:SER:O	3:G:43:ILE:HG13	2.16	0.45
1:E:39:TYR:HB2	1:E:42:ARG:HD3	1.97	0.45
1:E:321:GLU:C	1:E:323:LYS:H	2.25	0.45
2:B:242:PRO:HB3	4:H:93:VAL:C	2.41	0.45
1:E:36:SER:HA	5:E:601:FAD:N3A	2.31	0.45
2:F:116:ILE:HB	2:F:176:ALA:HB2	1.99	0.45
1:A:455:MET:HE2	1:A:455:MET:HB3	1.81	0.44
2:B:65:CYS:SG	2:B:76:ALA:HB3	2.57	0.44
1:E:304:ILE:HG13	1:E:313:TYR:CE2	2.51	0.44
2:F:154:CYS:HB2	2:F:210:CYS:HB2	1.98	0.44
1:A:198:VAL:O	1:A:456:GLY:HA2	2.18	0.44
1:E:42:ARG:HG3	2:F:62:CYS:SG	2.58	0.44
1:E:256:ASN:HB2	1:E:302:ASN:HB3	2.00	0.44
1:E:295:TRP:HE1	1:E:299:ARG:HE	1.64	0.44
1:A:182:GLN:OE1	1:A:184:ARG:NH1	2.44	0.44
2:B:110:ILE:O	2:B:114:GLU:HG3	2.18	0.44
1:E:230:GLN:HB2	1:E:358:MET:HE1	1.99	0.44
1:A:133:PHE:CE1	2:B:149:ILE:HG22	2.53	0.44
1:E:311:VAL:HG23	1:E:350:VAL:O	2.18	0.43
2:B:119:TYR:O	2:B:191:ARG:NH2	2.51	0.43
1:E:232:HIS:CE1	1:E:245:GLN:HG3	2.53	0.43
1:E:553:ALA:HA	1:E:563:LEU:HD23	2.00	0.43
1:A:386:HIS:CE1	1:A:390:ARG:HG3	2.53	0.43
4:H:63:GLY:O	4:H:67:LEU:HD23	2.18	0.43
1:A:308:ARG:HH12	1:A:339:VAL:HA	1.83	0.43
1:E:225:ASP:HB3	1:E:228:PHE:HD1	1.83	0.43
3:G:9:ARG:HE	3:G:9:ARG:HB2	1.48	0.43
1:A:392:GLY:O	1:A:393:SER:OG	2.32	0.43
1:E:42:ARG:HD2	1:E:42:ARG:N	2.33	0.43
1:E:42:ARG:NH2	2:F:63:GLY:O	2.52	0.43
2:B:116:ILE:HB	2:B:176:ALA:HB2	2.00	0.43
2:F:113:LEU:HD23	2:F:118:PRO:HG3	2.00	0.43
3:G:82:HIS:O	3:G:86:TRP:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:6:LEU:HD22	2:B:81:LEU:HB3	2.01	0.43
2:B:216:LYS:HE2	2:B:216:LYS:HA	2.00	0.43
1:E:270:PRO:HD2	1:E:281:TYR:CE2	2.49	0.43
2:F:229:VAL:HG12	2:F:233:LYS:HE3	2.00	0.43
1:E:217:LEU:HD11	1:E:517:CYS:SG	2.59	0.43
1:A:272:THR:H	1:A:282:MET:HG3	1.84	0.43
1:E:469:PRO:HB3	1:E:523:MET:HE1	2.00	0.42
2:F:117:LYS:HD2	2:F:119:TYR:OH	2.19	0.42
1:A:231:TYR:O	1:A:287:ARG:HD3	2.19	0.42
1:A:260:TYR:O	1:A:262:TYR:N	2.52	0.42
3:C:34:PRO:HA	3:C:37:TRP:HB3	2.01	0.42
4:D:63:GLY:O	4:D:67:LEU:HD23	2.19	0.42
1:E:363:THR:HB	1:E:367:CYS:HA	2.01	0.42
1:A:119:MET:HE2	1:A:391:LEU:HD23	2.00	0.42
2:F:39:ALA:O	2:F:43:ILE:HG13	2.20	0.42
1:E:148:GLN:OE1	1:E:148:GLN:N	2.41	0.42
3:G:33:VAL:HA	4:H:82:MET:HE3	2.01	0.42
1:A:256:ASN:HB2	1:A:302:ASN:O	2.20	0.42
1:A:327:ARG:HA	1:A:327:ARG:HD3	1.84	0.42
1:E:279:ASN:C	1:E:281:TYR:H	2.27	0.42
1:A:236:LEU:HD22	1:A:339:VAL:HG11	2.02	0.42
1:A:358:MET:HE3	1:A:389:ASN:HA	2.01	0.42
2:B:9:GLU:HG3	2:B:23:SER:HB3	2.02	0.42
3:G:34:PRO:HB3	3:G:75:THR:HA	2.02	0.42
1:A:46:VAL:HB	1:A:136:LEU:HD23	2.02	0.41
1:A:103:TRP:O	2:B:139:MET:HE1	2.20	0.41
4:D:10:GLU:N	4:D:11:PRO:HD2	2.35	0.41
2:F:155:TYR:CZ	2:F:169:GLY:HA3	2.54	0.41
1:A:228:PHE:O	1:A:358:MET:HB2	2.19	0.41
4:D:75:LEU:HD23	4:D:75:LEU:HA	1.90	0.41
1:A:44:HIS:CE1	1:A:204:ASN:HA	2.56	0.41
1:A:396:LEU:HD23	1:A:396:LEU:HA	1.81	0.41
4:H:112:LEU:HA	4:H:115:VAL:HG22	2.03	0.41
1:A:553:ALA:HA	1:A:563:LEU:HD23	2.03	0.41
1:E:399:LEU:HD11	5:E:601:FAD:H4'	2.02	0.41
1:E:498:PHE:CD2	2:F:103:VAL:HG21	2.56	0.41
2:B:116:ILE:H	2:B:116:ILE:HG13	1.70	0.41
4:D:93:VAL:C	2:F:242:PRO:HB3	2.45	0.41
2:B:39:ALA:O	2:B:43:ILE:HG13	2.21	0.41
2:F:241:LYS:HA	2:F:242:PRO:HD3	1.80	0.41
1:A:60:ASP:OD2	1:A:61:SER:N	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:TRP:CD1	1:A:125:TRP:N	2.89	0.41
1:A:230:GLN:HB2	1:A:358:MET:SD	2.61	0.41
3:C:98:VAL:HG23	3:C:98:VAL:O	2.21	0.41
1:E:217:LEU:HD23	1:E:217:LEU:HA	1.96	0.41
1:A:118:GLY:HA2	1:A:279:ASN:HD21	1.86	0.41
1:A:549:LYS:HB2	1:A:566:SER:O	2.21	0.41
2:B:4:LYS:O	2:B:29:PRO:HA	2.21	0.41
1:E:308:ARG:HD2	1:E:308:ARG:N	2.36	0.41
1:A:282:MET:HB3	1:A:283:GLU:OE1	2.21	0.40
1:A:308:ARG:NH1	1:A:339:VAL:HG12	2.37	0.40
1:A:363:THR:HB	1:A:367:CYS:HA	2.01	0.40
2:B:121:ILE:HB	2:B:187:GLY:HA3	2.02	0.40
4:D:28:ALA:O	4:D:32:ILE:HG13	2.20	0.40
1:E:42:ARG:HG2	2:F:64:SER:HB3	2.04	0.40
1:E:242:LEU:HB3	5:E:601:FAD:HM72	2.02	0.40
1:E:497:VAL:O	2:F:100:ARG:NH1	2.51	0.40
2:F:216:LYS:HA	2:F:216:LYS:HE2	2.03	0.40
3:G:124:LEU:HD21	4:H:33:LEU:HD23	2.02	0.40
1:A:232:HIS:H	1:A:355:HIS:HB2	1.86	0.40
2:B:57:CYS:HB3	2:B:62:CYS:HB3	2.03	0.40
2:B:241:LYS:HA	2:B:242:PRO:HD3	1.88	0.40
1:E:125:TRP:CD1	1:E:125:TRP:N	2.88	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	575/585 (98%)	540 (94%)	31 (5%)	4 (1%)	18	55
1	E	583/585 (100%)	538 (92%)	37 (6%)	8 (1%)	9	39
2	B	241/243 (99%)	225 (93%)	13 (5%)	3 (1%)	10	43

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	241/243 (99%)	230 (95%)	8 (3%)	3 (1%)	10	43
3	C	128/130 (98%)	122 (95%)	6 (5%)	0	100	100
3	G	128/130 (98%)	121 (94%)	6 (5%)	1 (1%)	16	52
4	D	117/119 (98%)	110 (94%)	6 (5%)	1 (1%)	14	49
4	H	117/119 (98%)	111 (95%)	5 (4%)	1 (1%)	14	49
All	All	2130/2154 (99%)	1997 (94%)	112 (5%)	21 (1%)	12	47

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	PRO
1	E	579	VAL
1	A	261	ARG
2	B	242	PRO
1	E	329	PRO
2	F	242	PRO
1	A	128	ALA
1	A	234	THR
4	D	117	THR
3	G	18	LYS
4	H	117	THR
1	E	128	ALA
1	E	273	PRO
1	E	322	LYS
2	F	56	SER
2	B	56	SER
1	E	312	VAL
2	B	61	ILE
2	F	61	ILE
1	E	444	GLY
1	E	581	GLY

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	460/465 (99%)	454 (99%)	6 (1%)	61	71
1	E	465/465 (100%)	449 (97%)	16 (3%)	32	54
2	B	205/205 (100%)	195 (95%)	10 (5%)	22	45
2	F	205/205 (100%)	195 (95%)	10 (5%)	22	45
3	C	111/111 (100%)	108 (97%)	3 (3%)	39	60
3	G	111/111 (100%)	106 (96%)	5 (4%)	24	47
4	D	97/97 (100%)	92 (95%)	5 (5%)	21	43
4	H	97/97 (100%)	93 (96%)	4 (4%)	27	49
All	All	1751/1756 (100%)	1692 (97%)	59 (3%)	32	54

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ARG
1	A	122	GLU
1	A	200	ARG
1	A	332	CYS
1	A	351	ARG
1	A	400	VAL
2	B	19	THR
2	B	62	CYS
2	B	65	CYS
2	B	128	ASP
2	B	148	CYS
2	B	151	CYS
2	B	204	CYS
2	B	212	GLU
2	B	239	THR
2	B	243	ARG
3	C	50	LYS
3	C	84	LYS
3	C	103	MET
4	D	7	ARG
4	D	9	ASP
4	D	86	MET
4	D	117	THR
4	D	118	ILE
1	E	18	ARG
1	E	96	LEU
1	E	122	GLU

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Mol	Chain	Res	Type
1	E	239	SER
1	E	243	MET
1	E	244	THR
1	E	268	MET
1	E	274	LEU
1	E	287	ARG
1	E	288	ASP
1	E	311	VAL
1	E	324	LEU
1	E	326	GLU
1	E	332	CYS
1	E	351	ARG
1	E	537	GLU
2	F	19	THR
2	F	62	CYS
2	F	65	CYS
2	F	128	ASP
2	F	148	CYS
2	F	151	CYS
2	F	203	SER
2	F	204	CYS
2	F	212	GLU
2	F	243	ARG
3	G	9	ARG
3	G	50	LYS
3	G	84	LYS
3	G	95	ASN
3	G	103	MET
4	H	6	LYS
4	H	86	MET
4	H	117	THR
4	H	118	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	HIS
1	A	258	ASN
1	A	279	ASN
1	A	386	HIS
1	A	428	GLN
1	A	533	GLN

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Mol	Chain	Res	Type
2	B	196	ASN
2	B	225	GLN
3	C	64	GLN
1	E	155	HIS
1	E	204	ASN
1	E	230	GLN
1	E	232	HIS
1	E	264	GLN
1	E	389	ASN
1	E	428	GLN
1	E	442	GLN
2	F	225	GLN
2	F	226	GLN
4	H	83	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

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	FAD	E	601	-	58,58,58	0.32	0	85,89,89	0.34	0
8	SF4	F	303	2	0,12,12	-	-	-		
6	FES	B	301	-	0,4,4	-	-	-		
5	FAD	A	601	-	58,58,58	0.29	0	85,89,89	0.31	0
8	SF4	B	303	-	0,12,12	-	-	-		
7	F3S	B	302	-	0,9,9	-	-	-		
6	FES	F	301	-	0,4,4	-	-	-		
7	F3S	F	302	-	0,9,9	-	-	-		

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	FAD	E	601	-	-	16/34/50/50	0/6/6/6
8	SF4	F	303	2	-	-	0/6/5/5
6	FES	B	301	-	-	-	0/1/1/1
5	FAD	A	601	-	-	15/34/50/50	0/6/6/6
8	SF4	B	303	-	-	-	0/6/5/5
7	F3S	B	302	-	-	-	0/3/3/3
6	FES	F	301	-	-	-	0/1/1/1
7	F3S	F	302	-	-	-	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	601	FAD	N10-C1'-C2'-O2'
5	A	601	FAD	N10-C1'-C2'-C3'
5	A	601	FAD	C2'-C3'-C4'-O4'
5	A	601	FAD	C2'-C3'-C4'-C5'
5	A	601	FAD	O3'-C3'-C4'-O4'
5	A	601	FAD	O3'-C3'-C4'-C5'
5	A	601	FAD	C5'-O5'-P-O1P
5	A	601	FAD	C5'-O5'-P-O2P
5	A	601	FAD	C5'-O5'-P-O3P
5	E	601	FAD	O4'-C4'-C5'-O5'

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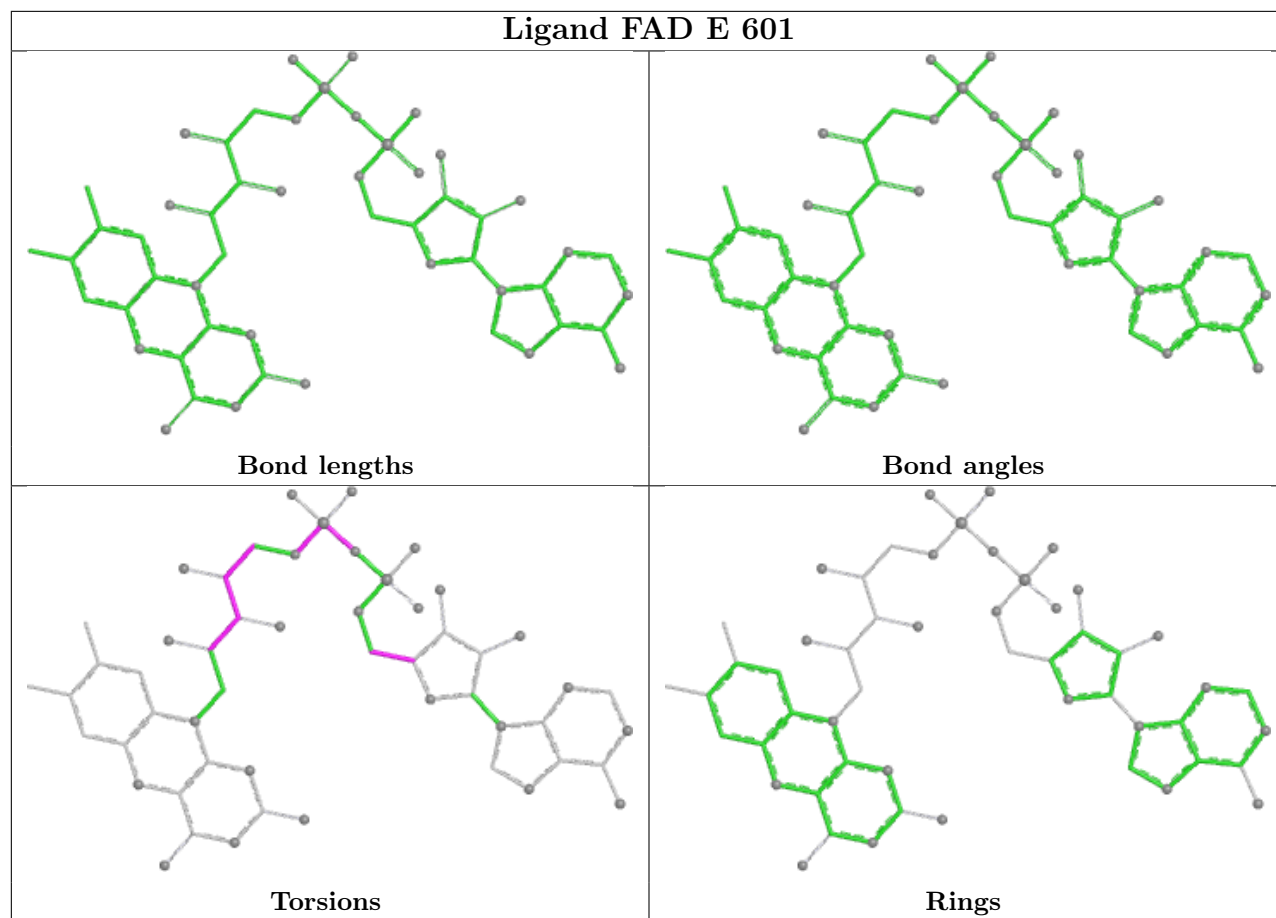
Mol	Chain	Res	Type	Atoms
5	E	601	FAD	C5'-O5'-P-O1P
5	E	601	FAD	C5'-O5'-P-O2P
5	E	601	FAD	C5'-O5'-P-O3P
5	E	601	FAD	PA-O3P-P-O5'
5	E	601	FAD	C2'-C3'-C4'-O4'
5	E	601	FAD	C3'-C4'-C5'-O5'
5	E	601	FAD	O3'-C3'-C4'-O4'
5	A	601	FAD	PA-O3P-P-O5'
5	E	601	FAD	C2'-C3'-C4'-C5'
5	A	601	FAD	O4'-C4'-C5'-O5'
5	E	601	FAD	O2'-C2'-C3'-O3'
5	E	601	FAD	O3'-C3'-C4'-C5'
5	E	601	FAD	O2'-C2'-C3'-C4'
5	E	601	FAD	PA-O3P-P-O1P
5	E	601	FAD	C1'-C2'-C3'-O3'
5	A	601	FAD	O2'-C2'-C3'-C4'
5	A	601	FAD	PA-O3P-P-O1P
5	A	601	FAD	O2'-C2'-C3'-O3'
5	A	601	FAD	O4B-C4B-C5B-O5B
5	E	601	FAD	PA-O3P-P-O2P
5	E	601	FAD	O4B-C4B-C5B-O5B

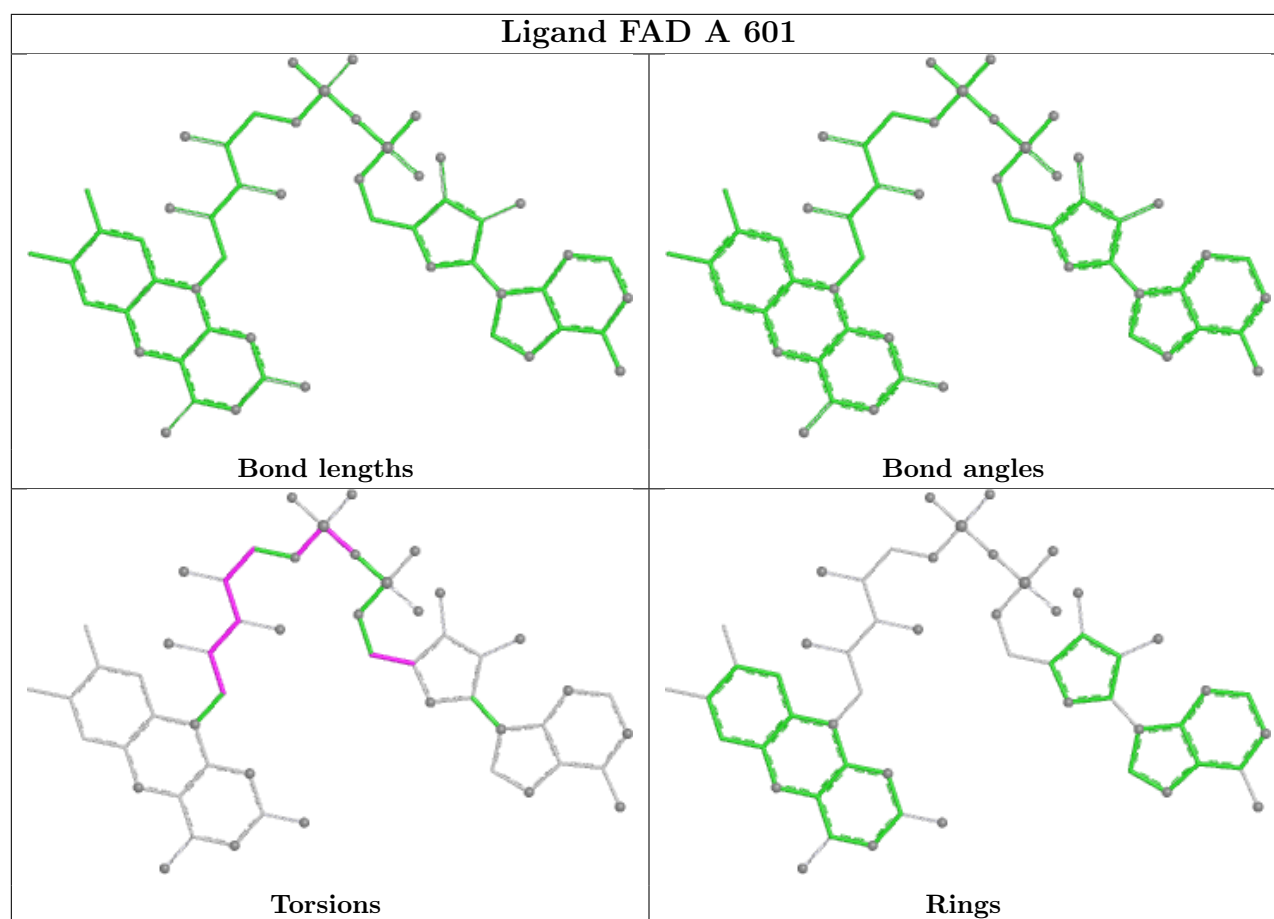
There are no ring outliers.

2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	601	FAD	10	0
5	A	601	FAD	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	577/585 (98%)	0.30	27 (4%) 36 32	155, 242, 339, 446	0
1	E	585/585 (100%)	0.20	19 (3%) 50 39	140, 249, 328, 385	0
2	B	243/243 (100%)	0.15	5 (2%) 63 49	125, 175, 237, 294	0
2	F	243/243 (100%)	0.21	5 (2%) 63 49	125, 179, 248, 287	0
3	C	130/130 (100%)	-0.09	2 (1%) 72 56	103, 157, 240, 295	0
3	G	130/130 (100%)	0.02	1 (0%) 82 68	92, 167, 271, 403	0
4	D	119/119 (100%)	-0.09	0 100 100	101, 146, 221, 276	0
4	H	119/119 (100%)	0.03	0 100 100	93, 141, 206, 279	0
All	All	2146/2154 (99%)	0.17	59 (2%) 56 43	92, 209, 318, 446	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	47	ALA	6.7
1	A	402	PHE	5.5
1	A	207	ILE	4.8
1	E	402	PHE	4.4
1	A	176	MET	3.8
1	E	135	MET	3.7
1	A	229	VAL	3.5
2	B	170	PRO	3.5
2	F	95	ASN	3.3
1	E	213	MET	3.2
1	E	498	PHE	3.2
1	A	499	ASN	3.1
1	E	338	TYR	3.1
1	A	338	TYR	3.1
1	A	436	LEU	3.0
1	E	383	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
2	F	17	VAL	2.9
1	A	65	HIS	2.9
1	A	348	ILE	2.8
1	E	117	GLY	2.8
2	F	149	ILE	2.7
1	A	385	LEU	2.7
2	B	147	GLY	2.7
1	E	199	TYR	2.7
3	C	49	LEU	2.7
1	E	337	ALA	2.7
1	E	229	VAL	2.6
1	E	236	LEU	2.6
1	E	195	ALA	2.6
2	F	151	CYS	2.5
1	A	497	VAL	2.5
1	E	333	GLU	2.4
2	B	146	SER	2.4
1	E	41	MET	2.4
1	A	479	LEU	2.4
1	A	382	SER	2.3
1	E	526	LYS	2.3
1	E	521	SER	2.3
2	B	151	CYS	2.2
1	A	33	ALA	2.2
1	E	411	THR	2.2
1	A	319	LEU	2.2
1	A	515	ALA	2.2
1	A	213	MET	2.2
1	A	359	GLY	2.1
1	A	32	ILE	2.1
1	A	460	GLU	2.1
2	B	145	PHE	2.1
1	A	180	LEU	2.1
2	F	38	ASP	2.1
1	E	81	VAL	2.1
1	A	316	LEU	2.1
3	C	22	TYR	2.1
1	A	102	PRO	2.1
3	G	79	ALA	2.0
1	E	385	LEU	2.0
1	A	383	VAL	2.0
1	A	128	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	165	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

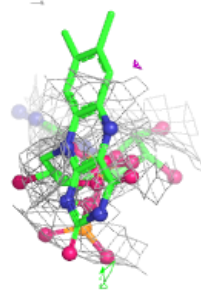
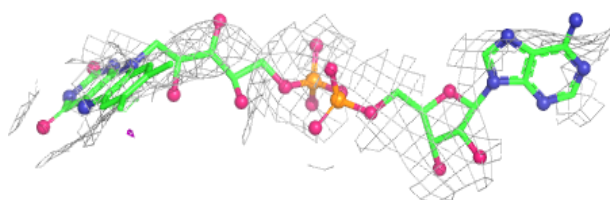
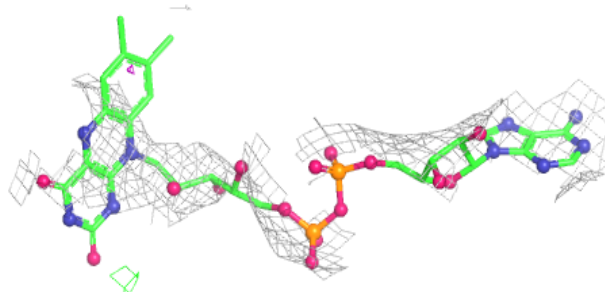
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	FAD	A	601	53/53	0.90	0.12	185,237,298,299	0
5	FAD	E	601	53/53	0.91	0.14	188,233,295,298	0
7	F3S	B	302	7/7	0.96	0.16	120,157,185,189	0
6	FES	F	301	4/4	0.97	0.11	142,149,154,158	0
8	SF4	F	303	8/8	0.98	0.09	127,147,160,162	0
7	F3S	F	302	7/7	0.99	0.09	101,175,201,217	0
8	SF4	B	303	8/8	0.99	0.09	131,137,151,162	0
6	FES	B	301	4/4	0.99	0.09	146,147,156,162	0

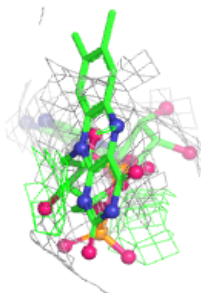
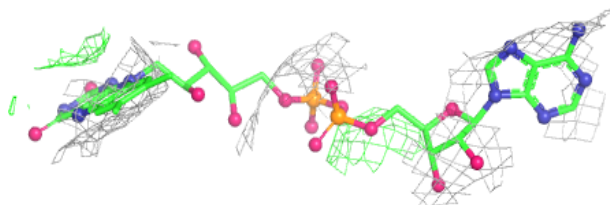
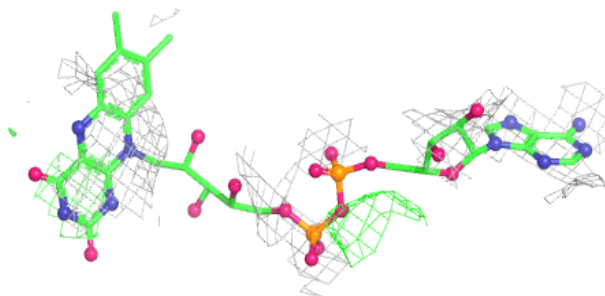
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around FAD A 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around FAD E 601:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.