



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 07:22 pm GMT

PDB ID : 3F90
Title : Desulfovibrio desulfuricans (ATCC 29577) semiquinone flavodoxin
Authors : Guelker, M.; Shamoo, Y.
Deposited on : 2008-11-13
Resolution : 2.50 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

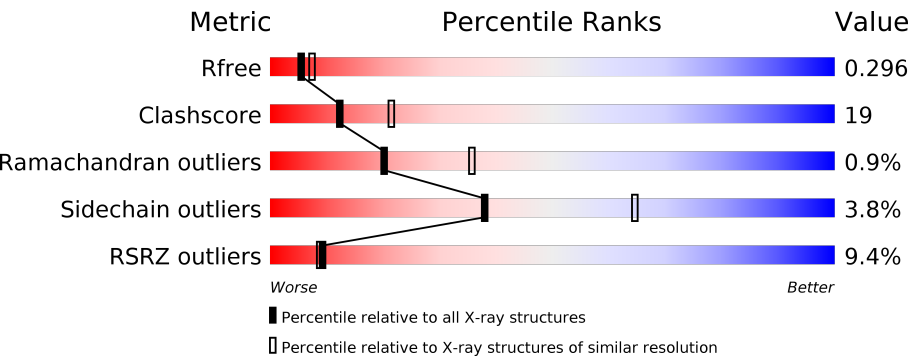
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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	148	<div><div>4%</div><div><div></div><div>72%</div><div>23%</div><div>5%</div></div><div></div></div>
1	B	148	<div><div>5%</div><div><div></div><div>70%</div><div>26%</div><div></div></div><div></div></div>
1	D	148	<div><div>17%</div><div><div></div><div>62%</div><div>35%</div><div></div></div><div></div></div>
1	E	148	<div><div>3%</div><div><div></div><div>68%</div><div>28%</div><div></div></div><div></div></div>
1	F	148	<div><div>12%</div><div><div></div><div>57%</div><div>40%</div><div></div></div><div></div></div>
1	G	148	<div><div>3%</div><div><div></div><div>68%</div><div>27%</div><div></div></div><div></div></div>

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Mol	Chain	Length	Quality of chain
1	H	148	<div><div></div><div>9%</div><div>61%</div><div>38%</div><div></div><div>.</div></div>
1	I	148	<div><div></div><div>20%</div><div>61%</div><div>36%</div><div></div><div>..</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9221 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 Flavodoxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	147	Total	C	N	O	S	0	0	0
			1093	683	175	230	5			
1	B	147	Total	C	N	O	S	0	0	0
			1093	683	175	230	5			
1	D	147	Total	C	N	O	S	0	0	0
			1093	683	175	230	5			
1	E	147	Total	C	N	O	S	0	0	0
			1093	683	175	230	5			
1	F	147	Total	C	N	O	S	0	0	0
			1093	683	175	230	5			
1	H	147	Total	C	N	O	S	0	0	0
			1093	683	175	230	5			
1	I	147	Total	C	N	O	S	0	0	0
			1093	683	175	230	5			
1	G	147	Total	C	N	O	S	0	0	0
			1093	683	175	230	5			

There are 16 discrepancies between the modelled and reference sequences:

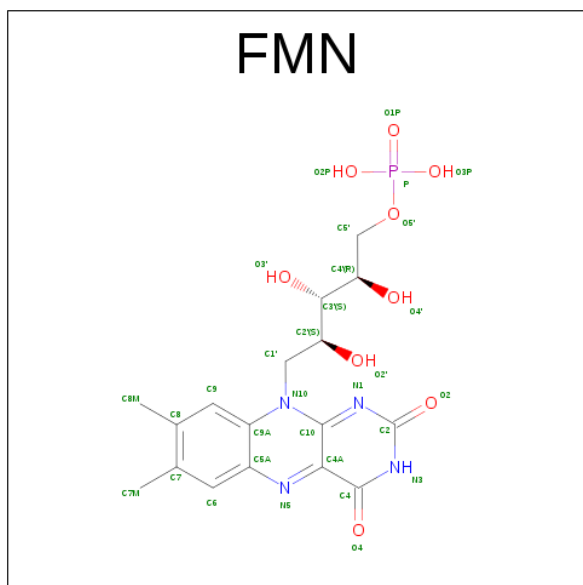
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP P26492
A	79	ASP	ASN	ENGINEERED	UNP P26492
B	1	MET	-	EXPRESSION TAG	UNP P26492
B	79	ASP	ASN	ENGINEERED	UNP P26492
D	1	MET	-	EXPRESSION TAG	UNP P26492
D	79	ASP	ASN	ENGINEERED	UNP P26492
E	1	MET	-	EXPRESSION TAG	UNP P26492
E	79	ASP	ASN	ENGINEERED	UNP P26492
F	1	MET	-	EXPRESSION TAG	UNP P26492
F	79	ASP	ASN	ENGINEERED	UNP P26492
H	1	MET	-	EXPRESSION TAG	UNP P26492
H	79	ASP	ASN	ENGINEERED	UNP P26492
I	1	MET	-	EXPRESSION TAG	UNP P26492

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Chain	Residue	Modelled	Actual	Comment	Reference
I	79	ASP	ASN	ENGINEERED	UNP P26492
G	1	MET	-	EXPRESSION TAG	UNP P26492
G	79	ASP	ASN	ENGINEERED	UNP P26492

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	D	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	E	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	F	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	H	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	I	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	G	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 3 is water.

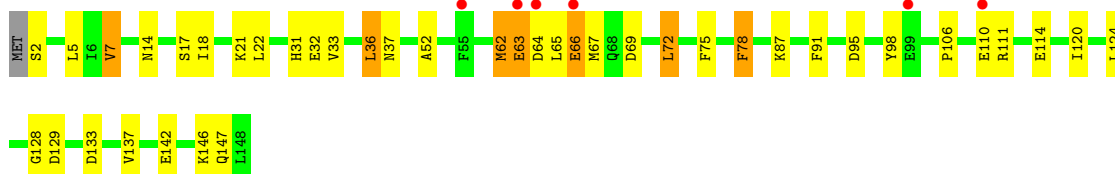
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	30	Total 30	O 30	0	0
3	B	33	Total 33	O 33	0	0
3	D	23	Total 23	O 23	0	0
3	E	35	Total 35	O 35	0	0
3	F	31	Total 31	O 31	0	0
3	H	31	Total 31	O 31	0	0
3	I	14	Total 14	O 14	0	0
3	G	32	Total 32	O 32	0	0

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 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.

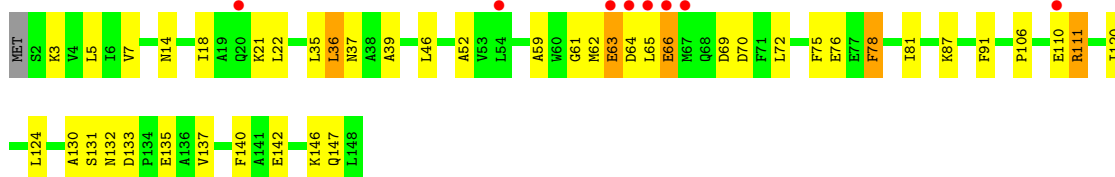
• Molecule 1: Flavodoxin

Chain A: 



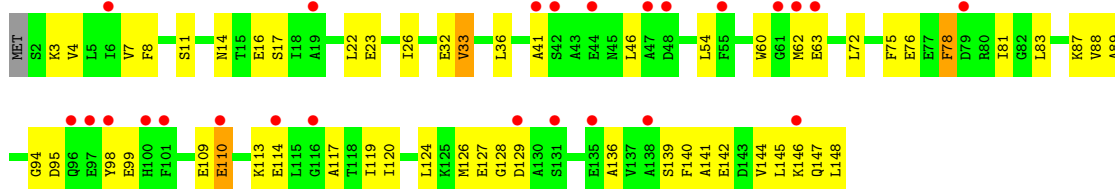
• Molecule 1: Flavodoxin

Chain B: 



• Molecule 1: Flavodoxin

Chain D: 



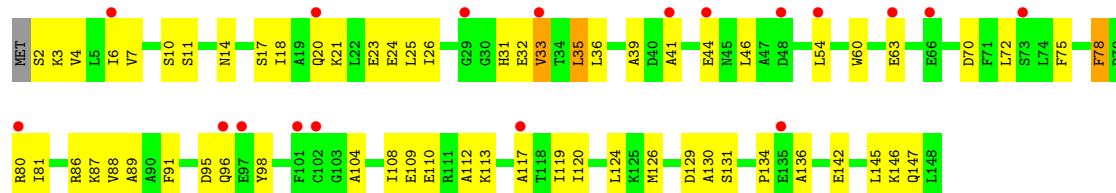
• Molecule 1: Flavodoxin

Chain E: 

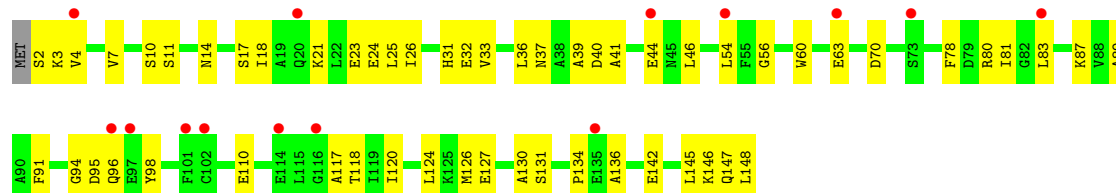




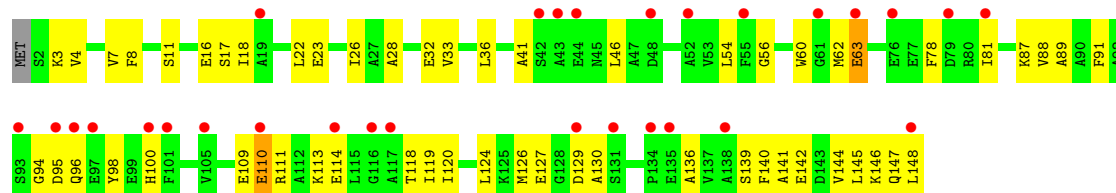
• Molecule 1: Flavodoxin



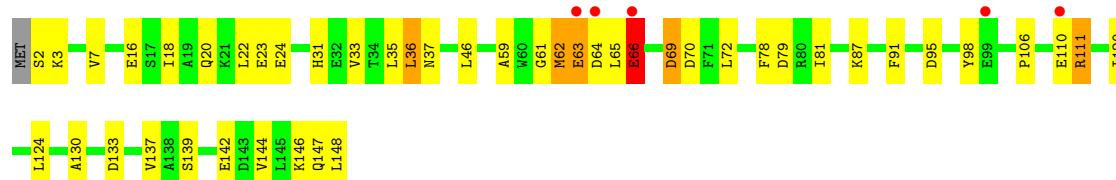
• Molecule 1: Flavodoxin



• Molecule 1: Flavodoxin



• Molecule 1: Flavodoxin



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	94.89Å 94.89Å 238.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.89 – 2.50 19.89 – 2.50	Depositor EDS
% Data completeness (in resolution range)	91.8 (19.89-2.50) 91.8 (19.89-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.79 (at 2.50Å)	Xtriage
Refinement program	CNS, PHENIX (phenix.refine)	Depositor
R, R_{free}	0.254 , 0.288 0.266 , 0.296	Depositor DCC
R_{free} test set	4042 reflections (10.05%)	DCC
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l	Xtriage
Reported twinning fraction	0.027 for -h,-k,l	Depositor
Outliers	3 of 40270 reflections (0.007%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9221	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 67.85 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.8084e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FMN

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	0.32	0/1108	0.49	0/1494
1	B	0.32	0/1108	0.48	0/1494
1	D	0.26	0/1108	0.41	0/1494
1	E	0.35	0/1108	0.50	0/1494
1	F	0.27	0/1108	0.40	0/1494
1	G	0.34	0/1108	0.49	0/1494
1	H	0.27	0/1108	0.41	0/1494
1	I	0.26	0/1108	0.41	0/1494
All	All	0.30	0/8864	0.45	0/11952

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

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	1093	0	1038	31	0
1	B	1093	0	1038	38	7
1	D	1093	0	1038	48	0
1	E	1093	0	1038	48	7
1	F	1093	0	1038	46	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	1093	0	1038	37	6
1	H	1093	0	1038	38	0
1	I	1093	0	1038	49	0
2	A	31	0	19	0	0
2	B	31	0	19	1	0
2	D	31	0	19	2	0
2	E	31	0	19	1	0
2	F	31	0	19	0	0
2	G	31	0	19	1	0
2	H	31	0	19	0	0
2	I	31	0	19	1	0
3	A	30	0	0	1	0
3	B	33	0	0	10	3
3	D	23	0	0	1	0
3	E	35	0	0	12	6
3	F	31	0	0	1	0
3	G	32	0	0	10	4
3	H	31	0	0	4	0
3	I	14	0	0	5	0
All	All	9221	0	8456	330	17

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 19.

All (330) 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:B:130:ALA:C	3:B:172:HOH:O	1.83	1.13
1:G:16:GLU:HG2	3:G:194:HOH:O	1.57	1.04
1:B:131:SER:C	3:B:172:HOH:O	1.95	1.03
1:B:132:ASN:N	3:B:172:HOH:O	1.90	1.03
1:H:40:ASP:OD2	3:H:221:HOH:O	1.82	0.97
1:G:20:GLN:CD	3:G:194:HOH:O	2.02	0.95
1:H:36:LEU:HD21	1:H:41:ALA:HB2	1.50	0.93
1:B:130:ALA:O	3:B:172:HOH:O	1.82	0.93
1:B:131:SER:N	3:B:172:HOH:O	1.96	0.92
1:B:131:SER:CA	3:B:172:HOH:O	2.18	0.89
1:E:78:PHE:O	3:E:149:HOH:O	1.90	0.89
1:E:69:ASP:OD1	3:E:155:HOH:O	1.92	0.88
1:I:126:MET:HE1	1:I:136:ALA:HB3	1.57	0.86
1:B:142:GLU:HG2	1:B:146:LYS:HE3	1.58	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:124:LEU:HD11	1:H:126:MET:HE2	1.57	0.86
1:F:36:LEU:HD21	1:F:41:ALA:HB2	1.59	0.85
1:G:20:GLN:OE1	3:G:194:HOH:O	1.96	0.83
1:E:44:GLU:CD	3:E:160:HOH:O	2.16	0.83
1:F:124:LEU:HD11	1:F:126:MET:HE2	1.61	0.82
1:E:142:GLU:HG2	1:E:146:LYS:HE3	1.62	0.81
1:H:118:THR:OG1	3:H:213:HOH:O	1.97	0.81
1:A:14:ASN:HD22	1:A:128:GLY:C	1.84	0.80
1:E:44:GLU:OE1	3:E:160:HOH:O	2.01	0.78
1:I:36:LEU:HD21	1:I:41:ALA:HB2	1.66	0.78
1:F:124:LEU:HD21	1:F:126:MET:HE3	1.63	0.78
1:G:70:ASP:O	3:G:157:HOH:O	2.03	0.76
1:F:142:GLU:O	1:F:146:LYS:HG2	1.86	0.76
1:I:111:ARG:NH2	3:I:160:HOH:O	2.00	0.76
1:E:44:GLU:OE2	3:E:160:HOH:O	2.04	0.76
1:D:36:LEU:HD21	1:D:41:ALA:HB2	1.68	0.76
1:D:126:MET:HE2	1:D:136:ALA:HB3	1.68	0.74
1:F:46:LEU:HB3	1:F:81:ILE:HG23	1.70	0.74
1:G:142:GLU:HG2	1:G:146:LYS:HE3	1.70	0.74
1:H:4:VAL:HB	1:H:33:VAL:HG12	1.68	0.74
1:H:23:GLU:HG3	1:H:33:VAL:HG21	1.70	0.73
1:E:106:PRO:O	1:E:110:GLU:HG3	1.89	0.72
1:I:95:ASP:HB3	1:I:98:TYR:HD1	1.53	0.72
1:F:54:LEU:HD22	1:F:89:ALA:HB3	1.71	0.71
1:E:3:LYS:NZ	3:E:169:HOH:O	2.22	0.71
1:F:11:SER:HB2	1:F:60:TRP:CZ2	2.27	0.70
1:H:46:LEU:HB3	1:H:81:ILE:HG23	1.73	0.70
1:B:65:LEU:HD12	1:B:72:LEU:HD21	1.74	0.69
1:I:4:VAL:HB	1:I:33:VAL:HG12	1.73	0.69
1:H:126:MET:HE1	1:H:136:ALA:HB3	1.73	0.69
1:H:142:GLU:O	1:H:146:LYS:HG2	1.93	0.69
1:A:62:MET:O	1:A:64:ASP:N	2.25	0.68
1:E:142:GLU:OE2	3:E:173:HOH:O	2.11	0.68
1:H:95:ASP:HB3	1:H:98:TYR:HD1	1.59	0.68
1:F:4:VAL:HB	1:F:33:VAL:HG12	1.76	0.68
1:H:3:LYS:HG2	1:H:32:GLU:HB2	1.76	0.68
1:D:124:LEU:HD11	1:D:126:MET:HE3	1.75	0.67
1:I:126:MET:HE1	1:I:136:ALA:CB	2.25	0.66
1:A:120:ILE:HB	1:A:147:GLN:HG3	1.78	0.66
1:D:95:ASP:HB3	1:D:98:TYR:HD1	1.60	0.65
1:H:124:LEU:HD11	1:H:126:MET:CE	2.26	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:69:ASP:OD2	3:E:171:HOH:O	2.14	0.65
1:G:3:LYS:NZ	3:G:168:HOH:O	2.29	0.65
1:G:139:SER:OG	3:G:173:HOH:O	2.06	0.64
1:G:65:LEU:HD12	1:G:72:LEU:HD21	1.78	0.64
1:H:37:ASN:HB3	3:H:221:HOH:O	1.97	0.64
1:I:26:ILE:HA	1:I:145:LEU:HD21	1.79	0.64
1:D:99:GLU:HB2	3:D:190:HOH:O	1.96	0.64
1:E:42:SER:N	3:E:220:HOH:O	2.17	0.63
1:B:75:PHE:HA	1:B:78:PHE:CD1	2.33	0.63
1:F:3:LYS:HG2	1:F:32:GLU:HB2	1.81	0.63
1:I:95:ASP:HB3	1:I:98:TYR:CD1	2.32	0.63
1:D:4:VAL:HB	1:D:33:VAL:HG12	1.80	0.63
1:F:23:GLU:HG3	1:F:33:VAL:HG21	1.80	0.63
1:H:120:ILE:HB	1:H:147:GLN:HG3	1.81	0.63
1:I:23:GLU:HG3	1:I:33:VAL:HG21	1.81	0.62
1:E:46:LEU:HB3	1:E:81:ILE:HG23	1.81	0.62
1:F:2:SER:N	1:F:31:HIS:HD1	1.98	0.62
1:F:95:ASP:HB3	1:F:98:TYR:HD1	1.65	0.61
1:G:106:PRO:O	1:G:110:GLU:HG3	1.99	0.61
1:D:142:GLU:O	1:D:146:LYS:HG2	2.00	0.61
1:F:124:LEU:HD11	1:F:126:MET:CE	2.30	0.61
1:H:23:GLU:HG3	1:H:33:VAL:CG2	2.30	0.61
1:D:46:LEU:HB3	1:D:81:ILE:HG23	1.82	0.61
1:A:106:PRO:O	1:A:110:GLU:HG3	2.00	0.61
1:A:21:LYS:HG2	3:A:165:HOH:O	1.99	0.61
1:B:3:LYS:NZ	3:B:154:HOH:O	2.33	0.61
1:H:124:LEU:HD21	1:H:126:MET:HE3	1.81	0.61
1:E:2:SER:N	1:E:31:HIS:HD1	1.99	0.61
1:H:21:LYS:O	1:H:25:LEU:HG	2.00	0.60
1:A:32:GLU:CD	1:F:20:GLN:HE21	2.05	0.60
1:G:20:GLN:NE2	3:G:194:HOH:O	2.30	0.60
1:E:145:LEU:O	3:E:162:HOH:O	2.17	0.60
1:G:69:ASP:OD1	3:G:150:HOH:O	2.17	0.60
1:I:120:ILE:HB	1:I:147:GLN:HG3	1.82	0.60
1:A:142:GLU:HG2	1:A:146:LYS:HE3	1.84	0.59
1:D:126:MET:HE2	1:D:136:ALA:CB	2.32	0.59
1:I:124:LEU:HD11	1:I:126:MET:HE2	1.83	0.59
1:E:14:ASN:HD22	1:E:129:ASP:HB3	1.67	0.58
1:D:95:ASP:HB3	1:D:98:TYR:CD1	2.37	0.58
1:A:14:ASN:HD22	1:A:129:ASP:N	2.01	0.58
1:H:7:VAL:HG13	1:H:36:LEU:HD22	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:46:LEU:HB3	1:G:81:ILE:HG23	1.85	0.58
1:F:126:MET:HE1	1:F:136:ALA:HB3	1.86	0.57
1:A:36:LEU:HD22	1:A:37:ASN:O	2.03	0.57
1:F:21:LYS:HD2	1:F:24:GLU:OE1	2.04	0.57
1:B:142:GLU:O	1:B:146:LYS:HG2	2.04	0.57
1:D:62:MET:CE	2:D:151:FMN:HM72	2.35	0.57
1:H:11:SER:HB2	1:H:60:TRP:CZ2	2.39	0.57
1:I:142:GLU:O	1:I:146:LYS:HG2	2.05	0.57
1:D:124:LEU:HD11	1:D:126:MET:CE	2.34	0.57
1:G:120:ILE:HB	1:G:147:GLN:HG3	1.87	0.56
1:F:7:VAL:HG13	1:F:36:LEU:HD22	1.87	0.56
1:I:87:LYS:NZ	3:I:182:HOH:O	2.29	0.56
1:D:120:ILE:HB	1:D:147:GLN:HG3	1.87	0.56
1:F:120:ILE:HB	1:F:147:GLN:HG3	1.87	0.56
1:I:94:GLY:O	1:I:127:GLU:HG2	2.05	0.56
1:A:65:LEU:O	1:A:66:GLU:HB2	2.05	0.56
1:B:106:PRO:O	1:B:110:GLU:HG3	2.06	0.56
1:H:87:LYS:HB3	1:H:120:ILE:HD13	1.87	0.56
1:H:96:GLN:NE2	3:H:170:HOH:O	2.37	0.56
1:I:54:LEU:HD22	1:I:89:ALA:HB3	1.86	0.56
1:B:18:ILE:O	1:B:22:LEU:HG	2.06	0.56
1:D:26:ILE:HA	1:D:145:LEU:HD21	1.87	0.56
1:E:107:ALA:HA	1:E:110:GLU:OE1	2.05	0.55
1:E:113:LYS:NZ	3:E:153:HOH:O	2.35	0.55
1:E:60:TRP:O	1:E:66:GLU:HA	2.06	0.55
1:F:23:GLU:HG3	1:F:33:VAL:CG2	2.36	0.55
1:D:36:LEU:CD2	1:D:41:ALA:HB2	2.36	0.55
1:D:3:LYS:HG2	1:D:32:GLU:HB2	1.88	0.55
1:F:36:LEU:CD2	1:F:41:ALA:HB2	2.35	0.55
1:H:2:SER:N	1:H:31:HIS:HD1	2.05	0.55
1:A:95:ASP:HB3	1:A:98:TYR:HD1	1.71	0.54
1:I:36:LEU:CD2	1:I:41:ALA:HB2	2.36	0.54
1:A:18:ILE:O	1:A:22:LEU:HG	2.07	0.54
1:H:26:ILE:HA	1:H:145:LEU:HD21	1.88	0.54
1:A:52:ALA:HB2	1:A:87:LYS:HB2	1.90	0.54
1:G:36:LEU:HD22	1:G:37:ASN:O	2.08	0.53
1:G:111:ARG:HG2	1:G:111:ARG:HH11	1.74	0.53
1:I:11:SER:HB2	1:I:60:TRP:CZ2	2.43	0.53
1:I:141:ALA:O	1:I:145:LEU:HG	2.08	0.53
1:F:44:GLU:HB2	1:F:80:ARG:NH2	2.24	0.53
1:A:65:LEU:O	1:A:66:GLU:CB	2.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:54:LEU:CD2	1:F:89:ALA:HB3	2.37	0.53
1:A:32:GLU:HG3	1:F:20:GLN:HG2	1.91	0.53
1:F:26:ILE:HA	1:F:145:LEU:HD21	1.91	0.53
1:E:61:GLY:HA2	1:E:66:GLU:HG3	1.90	0.53
1:E:16:GLU:O	1:E:20:GLN:HG3	2.09	0.52
1:B:142:GLU:CG	1:B:146:LYS:HE3	2.36	0.52
1:F:87:LYS:HB3	1:F:120:ILE:HD13	1.92	0.52
1:H:44:GLU:HB2	1:H:80:ARG:NH2	2.25	0.52
1:A:95:ASP:HB3	1:A:98:TYR:CD1	2.44	0.52
1:B:46:LEU:HB3	1:B:81:ILE:HG23	1.92	0.52
1:I:23:GLU:HG3	1:I:33:VAL:CG2	2.40	0.52
1:H:36:LEU:CD2	1:H:41:ALA:HB2	2.31	0.51
1:G:3:LYS:CE	3:G:168:HOH:O	2.57	0.51
1:I:62:MET:CE	2:I:155:FMN:HM72	2.41	0.51
1:I:63:GLU:O	1:I:100:HIS:NE2	2.39	0.51
1:D:139:SER:O	1:D:142:GLU:HB3	2.10	0.51
1:I:139:SER:O	1:I:142:GLU:HB3	2.11	0.51
1:A:133:ASP:O	1:A:137:VAL:HG23	2.10	0.51
1:D:62:MET:HE2	2:D:151:FMN:HM72	1.92	0.51
1:G:24:GLU:OE2	3:G:155:HOH:O	2.19	0.51
1:B:52:ALA:HB2	1:B:87:LYS:HB2	1.92	0.51
1:G:111:ARG:HG2	1:G:111:ARG:NH1	2.26	0.51
1:G:59:ALA:HB3	2:G:156:FMN:C2	2.41	0.51
1:B:36:LEU:HD22	1:B:37:ASN:O	2.10	0.50
1:E:14:ASN:ND2	1:E:129:ASP:HB3	2.26	0.50
1:B:59:ALA:HB3	2:B:150:FMN:C2	2.41	0.50
1:D:110:GLU:O	1:D:114:GLU:HG3	2.12	0.50
1:I:3:LYS:HG2	1:I:32:GLU:HB2	1.93	0.50
1:I:28:ALA:HB3	3:I:152:HOH:O	2.11	0.50
1:A:67:MET:SD	1:A:72:LEU:HD23	2.52	0.50
1:E:142:GLU:CG	1:E:146:LYS:HE3	2.36	0.50
1:D:94:GLY:O	1:D:127:GLU:HG2	2.12	0.50
1:E:2:SER:N	3:E:158:HOH:O	2.44	0.50
1:B:87:LYS:HB3	1:B:120:ILE:HG21	1.93	0.50
1:F:18:ILE:HA	1:F:130:ALA:HB2	1.94	0.50
1:D:109:GLU:O	1:D:113:LYS:HG3	2.12	0.49
1:D:23:GLU:HG3	1:D:33:VAL:HG21	1.94	0.49
1:E:142:GLU:O	1:E:146:LYS:HG2	2.12	0.49
1:H:94:GLY:O	1:H:127:GLU:HG2	2.12	0.49
1:B:14:ASN:O	1:B:18:ILE:HG13	2.12	0.49
1:B:61:GLY:HA2	1:B:66:GLU:HG3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:PHE:HA	1:A:78:PHE:CD1	2.48	0.48
1:I:18:ILE:HA	1:I:130:ALA:HB2	1.95	0.48
1:A:33:VAL:O	1:F:23:GLU:OE2	2.31	0.48
1:A:14:ASN:O	1:A:18:ILE:HG13	2.13	0.48
1:E:120:ILE:HB	1:E:147:GLN:HG3	1.95	0.48
1:G:36:LEU:HD22	1:G:37:ASN:N	2.28	0.48
1:B:72:LEU:O	1:B:76:GLU:HG3	2.13	0.48
1:A:91:PHE:CB	1:A:124:LEU:HB3	2.43	0.48
1:E:133:ASP:O	1:E:137:VAL:HG23	2.13	0.48
1:H:131:SER:O	1:H:134:PRO:HD3	2.14	0.48
1:E:124:LEU:HB2	1:E:140:PHE:CD1	2.48	0.48
1:I:140:PHE:O	1:I:144:VAL:HG23	2.14	0.48
1:D:120:ILE:O	1:D:147:GLN:HG3	2.13	0.48
1:D:141:ALA:O	1:D:145:LEU:HG	2.13	0.48
1:F:131:SER:O	1:F:134:PRO:HD3	2.14	0.48
1:E:36:LEU:HD22	1:E:37:ASN:N	2.29	0.48
1:G:62:MET:HB3	1:G:63:GLU:H	1.49	0.48
1:E:22:LEU:O	1:E:26:ILE:HG13	2.14	0.47
1:G:18:ILE:HA	1:G:130:ALA:HB2	1.95	0.47
1:D:8:PHE:CZ	1:D:16:GLU:HG3	2.48	0.47
1:A:7:VAL:HG13	1:A:36:LEU:HB3	1.97	0.47
1:E:36:LEU:HD22	1:E:37:ASN:O	2.15	0.47
1:F:104:ALA:O	1:F:108:ILE:HG13	2.14	0.47
1:F:39:ALA:HB2	1:F:70:ASP:CG	2.34	0.47
1:G:18:ILE:O	1:G:22:LEU:HG	2.14	0.47
1:I:56:GLY:HA2	1:I:91:PHE:O	2.15	0.47
1:H:54:LEU:HD22	1:H:89:ALA:HB3	1.95	0.47
1:D:145:LEU:HD23	1:D:148:LEU:HD12	1.95	0.47
1:I:111:ARG:NH1	3:I:160:HOH:O	2.47	0.47
1:F:91:PHE:CB	1:F:124:LEU:HB3	2.45	0.47
1:I:87:LYS:HD3	1:I:120:ILE:HG21	1.96	0.47
1:D:11:SER:HB2	1:D:60:TRP:CZ2	2.49	0.47
1:E:106:PRO:O	1:E:110:GLU:CG	2.61	0.47
1:E:75:PHE:HA	1:E:78:PHE:CD1	2.50	0.47
1:G:2:SER:N	1:G:31:HIS:HD1	2.13	0.47
1:B:111:ARG:HD3	3:B:165:HOH:O	2.14	0.46
1:F:14:ASN:O	1:F:17:SER:HB3	2.16	0.46
1:E:52:ALA:HB2	1:E:87:LYS:HB2	1.96	0.46
1:G:133:ASP:O	1:G:137:VAL:HG23	2.15	0.46
1:D:54:LEU:HD22	1:D:89:ALA:HB3	1.97	0.46
1:E:64:ASP:O	1:E:66:GLU:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:LEU:HG	1:A:7:VAL:HG22	1.98	0.46
1:B:120:ILE:HB	1:B:147:GLN:HG3	1.96	0.46
1:D:87:LYS:HB3	1:D:120:ILE:HG21	1.97	0.46
1:D:23:GLU:HG3	1:D:33:VAL:CG2	2.45	0.46
1:I:124:LEU:HD11	1:I:126:MET:CE	2.43	0.46
1:I:8:PHE:CZ	1:I:16:GLU:HG3	2.51	0.46
1:F:17:SER:OG	1:F:129:ASP:HB2	2.15	0.46
1:F:86:ARG:HH11	1:F:86:ARG:HG2	1.80	0.46
1:H:18:ILE:HA	1:H:130:ALA:HB2	1.97	0.46
1:G:142:GLU:O	1:G:146:LYS:HG2	2.16	0.46
1:F:109:GLU:O	1:F:113:LYS:HG3	2.16	0.46
1:G:144:VAL:O	1:G:148:LEU:HG	2.17	0.45
1:I:88:VAL:O	1:I:119:ILE:HA	2.17	0.45
1:G:35:LEU:HA	1:G:35:LEU:HD12	1.86	0.45
1:H:36:LEU:HD23	1:H:36:LEU:C	2.36	0.45
1:A:62:MET:HB3	1:A:63:GLU:H	1.47	0.45
1:E:111:ARG:HA	1:E:111:ARG:HD2	1.53	0.45
1:D:87:LYS:HB3	1:D:120:ILE:HD13	1.99	0.45
1:D:124:LEU:HB2	1:D:140:PHE:CD1	2.52	0.45
1:I:46:LEU:HB3	1:I:81:ILE:HG23	1.99	0.45
1:B:91:PHE:CB	1:B:124:LEU:HB3	2.47	0.45
1:D:140:PHE:O	1:D:144:VAL:HG23	2.16	0.45
1:E:17:SER:OG	1:E:129:ASP:HB2	2.16	0.45
1:B:62:MET:O	1:B:64:ASP:N	2.50	0.45
1:A:2:SER:N	1:A:31:HIS:HD1	2.15	0.44
1:E:14:ASN:HD22	1:E:129:ASP:CB	2.30	0.44
1:I:110:GLU:O	1:I:114:GLU:HG3	2.18	0.44
1:E:110:GLU:O	1:E:114:GLU:HG3	2.17	0.44
1:F:72:LEU:O	1:F:72:LEU:HD12	2.17	0.44
1:H:39:ALA:HB2	1:H:70:ASP:CG	2.38	0.44
1:A:111:ARG:HA	1:A:111:ARG:HD2	1.59	0.44
1:B:111:ARG:HG2	1:B:111:ARG:HH11	1.82	0.44
1:I:120:ILE:O	1:I:147:GLN:HG3	2.17	0.44
1:F:36:LEU:HD23	1:F:36:LEU:C	2.38	0.44
1:F:112:ALA:O	1:F:117:ALA:HB3	2.18	0.44
1:I:87:LYS:HB3	1:I:120:ILE:CG2	2.47	0.44
1:D:88:VAL:O	1:D:119:ILE:HA	2.18	0.43
1:D:7:VAL:HA	1:D:36:LEU:O	2.18	0.43
1:H:3:LYS:HE3	1:H:3:LYS:HB3	1.88	0.43
1:I:7:VAL:HA	1:I:36:LEU:O	2.18	0.43
1:B:63:GLU:HB3	3:B:163:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:GLU:HG3	1:E:64:ASP:OD2	2.19	0.43
1:B:124:LEU:HB2	1:B:140:PHE:CD1	2.54	0.43
1:D:87:LYS:HB3	1:D:120:ILE:CG2	2.49	0.43
1:F:96:GLN:NE2	3:F:163:HOH:O	2.50	0.43
1:D:14:ASN:HB3	1:D:128:GLY:O	2.18	0.43
1:F:87:LYS:HB3	1:F:120:ILE:HG21	2.00	0.43
1:H:95:ASP:HB3	1:H:98:TYR:CD1	2.46	0.43
1:I:109:GLU:O	1:I:113:LYS:HG3	2.19	0.43
1:B:133:ASP:O	1:B:137:VAL:HG23	2.19	0.43
1:B:65:LEU:HD12	1:B:72:LEU:CD2	2.45	0.43
1:E:62:MET:O	1:E:64:ASP:N	2.51	0.43
1:E:14:ASN:HD22	1:E:129:ASP:N	2.17	0.43
1:I:17:SER:OG	1:I:129:ASP:HB2	2.19	0.43
1:G:7:VAL:HG21	1:G:46:LEU:HD21	2.01	0.42
1:B:111:ARG:HG2	1:B:111:ARG:NH1	2.34	0.42
1:B:5:LEU:HG	1:B:7:VAL:HG23	2.01	0.42
1:G:91:PHE:CB	1:G:124:LEU:HB3	2.49	0.42
1:I:22:LEU:O	1:I:26:ILE:HG13	2.18	0.42
1:F:6:ILE:HB	1:F:35:LEU:CD1	2.50	0.42
1:B:65:LEU:HB2	1:B:72:LEU:CD2	2.49	0.42
1:D:22:LEU:O	1:D:26:ILE:HG13	2.18	0.42
1:D:78:PHE:HA	1:D:81:ILE:HG13	2.01	0.42
1:E:59:ALA:HB3	2:E:152:FMN:C2	2.50	0.42
1:G:23:GLU:HG3	1:G:33:VAL:HB	2.01	0.42
1:I:91:PHE:CB	1:I:124:LEU:HB3	2.49	0.42
1:H:56:GLY:HA2	1:H:91:PHE:O	2.20	0.42
1:I:87:LYS:HB3	1:I:120:ILE:HG21	2.00	0.42
1:D:17:SER:OG	1:D:129:ASP:HB2	2.19	0.42
1:E:64:ASP:C	1:E:66:GLU:N	2.73	0.42
1:F:95:ASP:HB3	1:F:98:TYR:CD1	2.51	0.42
1:H:21:LYS:HD2	1:H:24:GLU:OE1	2.20	0.42
1:E:87:LYS:HB3	1:E:120:ILE:HG21	2.01	0.42
1:I:145:LEU:HD23	1:I:148:LEU:HD12	2.01	0.42
1:D:87:LYS:HD3	1:D:120:ILE:HG21	2.02	0.41
1:I:96:GLN:C	1:I:98:TYR:H	2.24	0.41
1:D:36:LEU:C	1:D:36:LEU:HD23	2.40	0.41
1:A:110:GLU:O	1:A:114:GLU:HG3	2.19	0.41
1:B:21:LYS:HG2	3:B:157:HOH:O	2.21	0.41
1:F:88:VAL:O	1:F:119:ILE:HA	2.20	0.41
1:H:145:LEU:HD23	1:H:148:LEU:HD12	2.02	0.41
1:I:120:ILE:HB	1:I:147:GLN:CG	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:72:LEU:O	1:D:76:GLU:HG3	2.21	0.41
1:G:95:ASP:HB3	1:G:98:TYR:HD1	1.84	0.41
1:D:75:PHE:HA	1:D:78:PHE:CD1	2.56	0.41
1:G:36:LEU:CD2	1:G:37:ASN:N	2.84	0.41
1:H:83:LEU:HB3	1:H:117:ALA:HB2	2.03	0.41
1:A:14:ASN:O	1:A:17:SER:HB3	2.20	0.41
1:B:39:ALA:HB2	1:B:70:ASP:CG	2.41	0.41
1:G:62:MET:O	1:G:64:ASP:N	2.54	0.41
1:D:33:VAL:O	1:E:23:GLU:OE2	2.39	0.41
1:E:104:ALA:O	1:E:108:ILE:HG13	2.21	0.41
1:F:3:LYS:HB3	1:F:3:LYS:HE3	1.91	0.41
1:I:111:ARG:CZ	3:I:160:HOH:O	2.56	0.41
1:I:36:LEU:HD23	1:I:36:LEU:C	2.41	0.41
1:D:3:LYS:HB3	1:D:3:LYS:HE3	1.85	0.41
1:E:62:MET:HB3	1:E:63:GLU:H	1.47	0.41
1:G:87:LYS:HB3	1:G:120:ILE:HG21	2.03	0.41
1:D:120:ILE:HB	1:D:147:GLN:CG	2.50	0.40
1:G:61:GLY:HA2	1:G:66:GLU:OE1	2.22	0.40
1:I:8:PHE:CE1	1:I:16:GLU:HG3	2.56	0.40
1:H:14:ASN:O	1:H:17:SER:HB3	2.20	0.40
1:F:75:PHE:HA	1:F:78:PHE:CD1	2.56	0.40
1:A:52:ALA:CB	1:A:87:LYS:HB2	2.50	0.40
1:B:35:LEU:HA	1:B:35:LEU:HD12	1.87	0.40
1:A:62:MET:C	1:A:64:ASP:H	2.23	0.40
1:D:83:LEU:HB3	1:D:117:ALA:HB2	2.04	0.40
1:E:17:SER:HA	1:E:20:GLN:OE1	2.21	0.40
1:F:21:LYS:O	1:F:25:LEU:HG	2.22	0.40
1:G:111:ARG:HH11	1:G:111:ARG:CG	2.35	0.40
1:I:87:LYS:HA	1:I:118:THR:O	2.21	0.40

All (17) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:66:GLU:CG	3:E:174:HOH:O[4_555]	0.58	1.62
1:E:66:GLU:OE2	3:E:156:HOH:O[4_555]	0.63	1.57
1:G:66:GLU:OE2	3:B:162:HOH:O[4_565]	0.67	1.53
1:G:66:GLU:CD	3:B:162:HOH:O[4_565]	0.88	1.32
1:E:66:GLU:CD	3:E:156:HOH:O[4_555]	0.90	1.30
1:E:66:GLU:CB	3:E:174:HOH:O[4_555]	1.07	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:GLU:CD	3:G:152:HOH:O[4_455]	1.14	1.06
1:G:66:GLU:OE1	3:B:162:HOH:O[4_565]	1.56	0.64
1:B:66:GLU:OE2	3:G:152:HOH:O[4_455]	1.56	0.64
1:E:66:GLU:OE1	3:E:156:HOH:O[4_555]	1.60	0.60
1:E:66:GLU:OE1	1:E:79:ASP:OD2[4_555]	1.62	0.58
1:B:66:GLU:CG	3:G:152:HOH:O[4_455]	1.84	0.36
1:B:66:GLU:OE1	3:G:152:HOH:O[4_455]	1.85	0.35
1:B:66:GLU:OE2	1:G:79:ASP:OD2[4_455]	1.86	0.34
1:B:66:GLU:OE1	1:G:79:ASP:OD2[4_455]	1.97	0.23
1:E:66:GLU:CD	3:E:174:HOH:O[4_555]	1.98	0.22
1:B:66:GLU:CD	1:G:79:ASP:OD2[4_455]	2.12	0.08

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	145/148 (98%)	138 (95%)	4 (3%)	3 (2%)	8	13
1	B	145/148 (98%)	138 (95%)	5 (3%)	2 (1%)	13	23
1	D	145/148 (98%)	136 (94%)	9 (6%)	0	100	100
1	E	145/148 (98%)	137 (94%)	6 (4%)	2 (1%)	13	23
1	F	145/148 (98%)	137 (94%)	8 (6%)	0	100	100
1	G	145/148 (98%)	136 (94%)	5 (3%)	4 (3%)	6	8
1	H	145/148 (98%)	138 (95%)	7 (5%)	0	100	100
1	I	145/148 (98%)	133 (92%)	12 (8%)	0	100	100
All	All	1160/1184 (98%)	1093 (94%)	56 (5%)	11 (1%)	20	36

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	63	GLU

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Mol	Chain	Res	Type
1	E	69	ASP
1	A	69	ASP
1	B	63	GLU
1	E	63	GLU
1	G	66	GLU
1	G	69	ASP
1	A	62	MET
1	B	69	ASP
1	G	62	MET
1	G	63	GLU

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	110/111 (99%)	105 (96%)	5 (4%)	32	56
1	B	110/111 (99%)	106 (96%)	4 (4%)	40	67
1	D	110/111 (99%)	106 (96%)	4 (4%)	40	67
1	E	110/111 (99%)	107 (97%)	3 (3%)	50	77
1	F	110/111 (99%)	104 (94%)	6 (6%)	25	46
1	G	110/111 (99%)	106 (96%)	4 (4%)	40	67
1	H	110/111 (99%)	106 (96%)	4 (4%)	40	67
1	I	110/111 (99%)	107 (97%)	3 (3%)	50	77
All	All	880/888 (99%)	847 (96%)	33 (4%)	38	64

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	36	LEU
1	A	66	GLU
1	A	72	LEU
1	A	78	PHE

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Mol	Chain	Res	Type
1	B	36	LEU
1	B	66	GLU
1	B	78	PHE
1	B	111	ARG
1	D	33	VAL
1	D	63	GLU
1	D	78	PHE
1	D	110	GLU
1	E	17	SER
1	E	36	LEU
1	E	78	PHE
1	F	10	SER
1	F	33	VAL
1	F	35	LEU
1	F	63	GLU
1	F	78	PHE
1	F	110	GLU
1	H	10	SER
1	H	63	GLU
1	H	78	PHE
1	H	110	GLU
1	I	63	GLU
1	I	78	PHE
1	I	110	GLU
1	G	36	LEU
1	G	66	GLU
1	G	78	PHE
1	G	111	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	E	14	ASN
1	F	20	GLN

5.3.3 RNA ⓘ

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 carbohydrates 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	$\# Z > 2$
2	FMN	A	149	-	31,33,33	1.39	4 (12%)	38,50,50	1.53	4 (10%)
2	FMN	B	150	-	31,33,33	1.41	4 (12%)	38,50,50	1.50	4 (10%)
2	FMN	D	151	-	31,33,33	1.40	5 (16%)	38,50,50	1.46	5 (13%)
2	FMN	E	152	-	31,33,33	1.44	4 (12%)	38,50,50	1.89	7 (18%)
2	FMN	F	153	-	31,33,33	1.41	5 (16%)	38,50,50	1.44	4 (10%)
2	FMN	G	156	-	31,33,33	1.41	4 (12%)	38,50,50	1.57	4 (10%)
2	FMN	H	154	-	31,33,33	1.41	5 (16%)	38,50,50	1.41	4 (10%)
2	FMN	I	155	-	31,33,33	1.42	5 (16%)	38,50,50	1.44	4 (10%)

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
2	FMN	A	149	-	-	0/16/18/18	0/3/3/3
2	FMN	B	150	-	-	0/16/18/18	0/3/3/3
2	FMN	D	151	-	-	0/16/18/18	0/3/3/3
2	FMN	E	152	-	-	0/16/18/18	0/3/3/3
2	FMN	F	153	-	-	0/16/18/18	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FMN	G	156	-	-	0/16/18/18	0/3/3/3
2	FMN	H	154	-	-	0/16/18/18	0/3/3/3
2	FMN	I	155	-	-	0/16/18/18	0/3/3/3

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	150	FMN	C5A-N5	2.05	1.38	1.35
2	F	153	FMN	C1'-N10	2.08	1.50	1.48
2	D	151	FMN	C1'-N10	2.08	1.50	1.48
2	D	151	FMN	C5A-N5	2.08	1.38	1.35
2	H	154	FMN	C5A-N5	2.12	1.38	1.35
2	G	156	FMN	C5A-N5	2.13	1.38	1.35
2	H	154	FMN	C1'-N10	2.13	1.50	1.48
2	A	149	FMN	C5A-N5	2.16	1.38	1.35
2	F	153	FMN	C5A-N5	2.19	1.38	1.35
2	I	155	FMN	C1'-N10	2.22	1.50	1.48
2	I	155	FMN	C5A-N5	2.22	1.38	1.35
2	E	152	FMN	C5A-N5	2.53	1.39	1.35
2	H	154	FMN	C4-N3	2.92	1.38	1.33
2	B	150	FMN	C4-N3	3.02	1.38	1.33
2	E	152	FMN	C4-N3	3.03	1.38	1.33
2	F	153	FMN	C4-N3	3.05	1.38	1.33
2	A	149	FMN	C4-N3	3.13	1.38	1.33
2	D	151	FMN	C4-N3	3.20	1.38	1.33
2	G	156	FMN	C4-N3	3.22	1.38	1.33
2	I	155	FMN	C4-N3	3.28	1.39	1.33
2	I	155	FMN	C4A-N5	3.66	1.38	1.33
2	D	151	FMN	C4A-N5	3.69	1.38	1.33
2	G	156	FMN	C4A-N5	3.74	1.38	1.33
2	A	149	FMN	C4A-N5	3.77	1.38	1.33
2	F	153	FMN	C10-N1	3.78	1.38	1.33
2	B	150	FMN	C4A-N5	3.82	1.38	1.33
2	A	149	FMN	C10-N1	3.83	1.38	1.33
2	H	154	FMN	C4A-N5	3.87	1.38	1.33
2	E	152	FMN	C10-N1	3.88	1.38	1.33
2	H	154	FMN	C10-N1	3.89	1.38	1.33
2	G	156	FMN	C10-N1	3.92	1.38	1.33
2	F	153	FMN	C4A-N5	3.96	1.39	1.33
2	I	155	FMN	C10-N1	3.98	1.38	1.33
2	D	151	FMN	C10-N1	4.02	1.38	1.33
2	B	150	FMN	C10-N1	4.07	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	152	FMN	C4A-N5	4.11	1.39	1.33

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	152	FMN	C4'-C3'-C2'	-3.91	104.98	113.41
2	E	152	FMN	C4A-C4-N3	-3.25	118.85	123.48
2	E	152	FMN	O5'-C5'-C4'	-3.21	100.80	109.36
2	A	149	FMN	C4A-C4-N3	-3.03	119.17	123.48
2	B	150	FMN	C4A-C4-N3	-2.94	119.29	123.48
2	G	156	FMN	C4A-C4-N3	-2.91	119.35	123.48
2	F	153	FMN	C4A-C4-N3	-2.58	119.81	123.48
2	D	151	FMN	C4A-C4-N3	-2.52	119.90	123.48
2	I	155	FMN	C4A-C4-N3	-2.50	119.92	123.48
2	H	154	FMN	C4A-C4-N3	-2.45	119.99	123.48
2	E	152	FMN	C4-C4A-N5	2.22	121.11	118.68
2	D	151	FMN	C1'-N10-C9A	2.43	120.58	118.35
2	B	150	FMN	C4A-N5-C5A	2.97	119.90	116.76
2	B	150	FMN	C5A-C9A-N10	3.04	119.91	117.66
2	F	153	FMN	C4A-N5-C5A	3.05	119.98	116.76
2	E	152	FMN	C4A-N5-C5A	3.06	119.99	116.76
2	F	153	FMN	C5A-C9A-N10	3.07	119.94	117.66
2	G	156	FMN	C4A-N5-C5A	3.12	120.06	116.76
2	A	149	FMN	C4A-N5-C5A	3.14	120.08	116.76
2	H	154	FMN	C5A-C9A-N10	3.15	120.00	117.66
2	I	155	FMN	C4A-N5-C5A	3.16	120.10	116.76
2	H	154	FMN	C4A-N5-C5A	3.17	120.11	116.76
2	D	151	FMN	C4A-N5-C5A	3.22	120.16	116.76
2	D	151	FMN	C5A-C9A-N10	3.23	120.06	117.66
2	I	155	FMN	C5A-C9A-N10	3.25	120.07	117.66
2	A	149	FMN	C5A-C9A-N10	3.35	120.15	117.66
2	E	152	FMN	C5A-C9A-N10	3.41	120.19	117.66
2	G	156	FMN	C5A-C9A-N10	3.72	120.42	117.66
2	I	155	FMN	C4-N3-C2	5.71	120.16	115.16
2	H	154	FMN	C4-N3-C2	5.72	120.16	115.16
2	D	151	FMN	C4-N3-C2	5.78	120.22	115.16
2	F	153	FMN	C4-N3-C2	5.93	120.35	115.16
2	G	156	FMN	C4-N3-C2	6.14	120.53	115.16
2	A	149	FMN	C4-N3-C2	6.27	120.64	115.16
2	B	150	FMN	C4-N3-C2	6.59	120.93	115.16
2	E	152	FMN	C4-N3-C2	7.07	121.34	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	150	FMN	1	0
2	D	151	FMN	2	0
2	E	152	FMN	1	0
2	G	156	FMN	1	0
2	I	155	FMN	1	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	147/148 (99%)	0.29	6 (4%) 38 40	26, 38, 57, 88	0
1	B	147/148 (99%)	0.27	8 (5%) 26 27	27, 39, 56, 89	0
1	D	147/148 (99%)	1.14	25 (17%) 2 1	40, 53, 74, 96	0
1	E	147/148 (99%)	0.28	5 (3%) 46 48	27, 38, 56, 89	0
1	F	147/148 (99%)	0.84	18 (12%) 5 4	41, 52, 74, 96	0
1	G	147/148 (99%)	0.26	5 (3%) 46 48	28, 38, 56, 89	0
1	H	147/148 (99%)	0.82	14 (9%) 9 8	41, 52, 74, 96	0
1	I	147/148 (99%)	1.17	29 (19%) 1 1	40, 53, 74, 96	0
All	All	1176/1184 (99%)	0.63	110 (9%) 9 9	26, 47, 70, 96	0

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	63	GLU	12.6
1	G	63	GLU	10.3
1	B	63	GLU	8.7
1	A	63	GLU	6.2
1	D	63	GLU	5.1
1	D	42	SER	5.0
1	I	61	GLY	4.6
1	E	64	ASP	4.5
1	D	97	GLU	4.4
1	I	97	GLU	4.3
1	F	135	GLU	4.3
1	I	63	GLU	4.3
1	G	64	ASP	4.2
1	D	61	GLY	4.2
1	H	96	GLN	4.0
1	I	44	GLU	3.9

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Mol	Chain	Res	Type	RSRZ
1	I	135	GLU	3.9
1	I	42	SER	3.8
1	I	117	ALA	3.7
1	I	110	GLU	3.7
1	D	138	ALA	3.6
1	D	110	GLU	3.6
1	E	65	LEU	3.5
1	D	48	ASP	3.4
1	D	135	GLU	3.4
1	B	64	ASP	3.4
1	I	43	ALA	3.4
1	B	110	GLU	3.3
1	F	20	GLN	3.2
1	H	54	LEU	3.2
1	D	62	MET	3.1
1	B	66	GLU	3.1
1	H	20	GLN	3.0
1	I	19	ALA	3.0
1	H	102	CYS	3.0
1	I	138	ALA	3.0
1	F	63	GLU	3.0
1	D	129	ASP	3.0
1	F	96	GLN	3.0
1	I	96	GLN	2.9
1	H	63	GLU	2.9
1	D	47	ALA	2.9
1	A	99	GLU	2.8
1	F	6	ILE	2.8
1	F	54	LEU	2.8
1	I	116	GLY	2.8
1	H	4	VAL	2.7
1	H	135	GLU	2.7
1	D	44	GLU	2.7
1	D	96	GLN	2.7
1	D	116	GLY	2.7
1	F	97	GLU	2.7
1	I	93	SER	2.7
1	D	55	PHE	2.7
1	I	100	HIS	2.6
1	E	110	GLU	2.6
1	I	131	SER	2.6
1	A	64	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	66	GLU	2.6
1	F	44	GLU	2.6
1	I	79	ASP	2.6
1	F	80	ARG	2.5
1	D	131	SER	2.5
1	D	19	ALA	2.5
1	I	95	ASP	2.5
1	D	114	GLU	2.5
1	D	98	TYR	2.4
1	F	102	CYS	2.4
1	F	29	GLY	2.4
1	H	44	GLU	2.4
1	F	48	ASP	2.4
1	H	114	GLU	2.4
1	F	117	ALA	2.3
1	D	6	ILE	2.3
1	I	81	ILE	2.3
1	B	67	MET	2.3
1	B	65	LEU	2.3
1	G	99	GLU	2.3
1	A	110	GLU	2.3
1	H	83	LEU	2.3
1	I	48	ASP	2.3
1	F	101	PHE	2.3
1	I	52	ALA	2.3
1	H	73	SER	2.2
1	I	55	PHE	2.2
1	I	76	GLU	2.2
1	I	148	LEU	2.2
1	G	66	GLU	2.2
1	D	100	HIS	2.2
1	A	55	PHE	2.2
1	I	101	PHE	2.2
1	H	116	GLY	2.2
1	D	79	ASP	2.2
1	I	105	VAL	2.2
1	H	97	GLU	2.2
1	B	54	LEU	2.2
1	I	134	PRO	2.1
1	D	101	PHE	2.1
1	F	33	VAL	2.1
1	F	41	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	I	114	GLU	2.1
1	B	20	GLN	2.1
1	E	44	GLU	2.1
1	I	129	ASP	2.1
1	F	66	GLU	2.1
1	F	73	SER	2.1
1	D	41	ALA	2.0
1	H	101	PHE	2.0
1	G	110	GLU	2.0
1	D	146	LYS	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 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, 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
2	FMN	H	154	31/31	0.88	0.22	0.43	48,56,59,61	0
2	FMN	A	149	31/31	0.94	0.17	0.22	38,42,44,52	0
2	FMN	F	153	31/31	0.87	0.22	0.21	51,56,58,61	0
2	FMN	G	156	31/31	0.94	0.17	-0.03	36,40,44,51	0
2	FMN	E	152	31/31	0.95	0.18	-0.04	38,41,45,52	0
2	FMN	I	155	31/31	0.91	0.22	-0.13	56,65,68,69	0
2	FMN	B	150	31/31	0.96	0.15	-0.19	37,40,44,50	0
2	FMN	D	151	31/31	0.90	0.21	-0.40	55,64,68,69	0

6.5 Other polymers [i](#)

There are no such residues in this entry.