



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:53 PM UTC

PDB ID : 3QFA / pdb_00003qfa
Title : Crystal structure of the human thioredoxin reductase-thioredoxin complex
Authors : Fritz-Wolf, K.; Kehr, S.; Stumpf, M.; Rahlfs, S.; Becker, K.
Deposited on : 2011-01-21
Resolution : 2.20 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

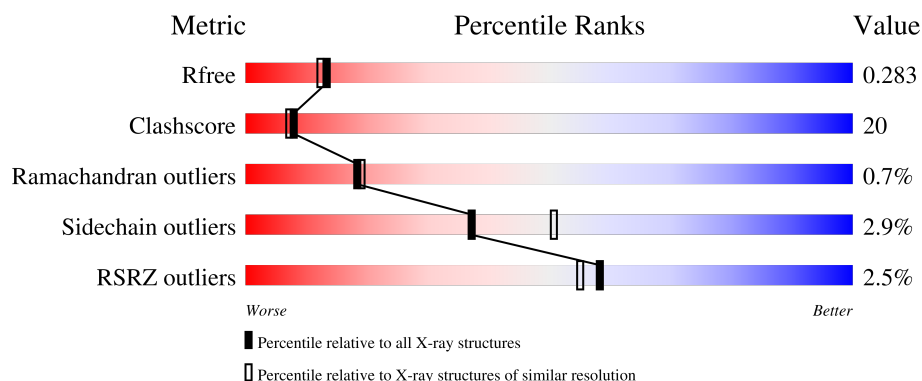
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	519	63% 29% • 5%
1	B	519	64% 29% • 5%
2	C	116	11% 48% 39% • 9%
2	D	116	4% 48% 40% • 9%

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
4	GOL	A	801	-	X	-	-
4	GOL	A	802	-	X	-	-
4	GOL	A	803	-	X	-	-
4	GOL	A	804	-	X	-	-
4	GOL	B	804	-	X	-	-
4	GOL	D	803	-	X	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9821 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 Thioredoxin reductase 1, cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	494	Total	C	N	O	S	0	0	0
			3808	2418	650	720	20			
1	B	494	Total	C	N	O	S	0	0	0
			3808	2418	650	720	20			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q16881
A	-18	GLY	-	expression tag	UNP Q16881
A	-17	SER	-	expression tag	UNP Q16881
A	-16	SER	-	expression tag	UNP Q16881
A	-15	HIS	-	expression tag	UNP Q16881
A	-14	HIS	-	expression tag	UNP Q16881
A	-13	HIS	-	expression tag	UNP Q16881
A	-12	HIS	-	expression tag	UNP Q16881
A	-11	HIS	-	expression tag	UNP Q16881
A	-10	HIS	-	expression tag	UNP Q16881
A	-9	SER	-	expression tag	UNP Q16881
A	-8	SER	-	expression tag	UNP Q16881
A	-7	GLY	-	expression tag	UNP Q16881
A	-6	LEU	-	expression tag	UNP Q16881
A	-5	VAL	-	expression tag	UNP Q16881
A	-4	PRO	-	expression tag	UNP Q16881
A	-3	ARG	-	expression tag	UNP Q16881
A	-2	GLY	-	expression tag	UNP Q16881
A	-1	SER	-	expression tag	UNP Q16881
A	0	HIS	-	expression tag	UNP Q16881
A	497	SER	CYS	engineered mutation	UNP Q16881
A	498	CYS	U	SEE REMARK 999	UNP Q16881
B	-19	MET	-	expression tag	UNP Q16881
B	-18	GLY	-	expression tag	UNP Q16881
B	-17	SER	-	expression tag	UNP Q16881

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	SER	-	expression tag	UNP Q16881
B	-15	HIS	-	expression tag	UNP Q16881
B	-14	HIS	-	expression tag	UNP Q16881
B	-13	HIS	-	expression tag	UNP Q16881
B	-12	HIS	-	expression tag	UNP Q16881
B	-11	HIS	-	expression tag	UNP Q16881
B	-10	HIS	-	expression tag	UNP Q16881
B	-9	SER	-	expression tag	UNP Q16881
B	-8	SER	-	expression tag	UNP Q16881
B	-7	GLY	-	expression tag	UNP Q16881
B	-6	LEU	-	expression tag	UNP Q16881
B	-5	VAL	-	expression tag	UNP Q16881
B	-4	PRO	-	expression tag	UNP Q16881
B	-3	ARG	-	expression tag	UNP Q16881
B	-2	GLY	-	expression tag	UNP Q16881
B	-1	SER	-	expression tag	UNP Q16881
B	0	HIS	-	expression tag	UNP Q16881
B	497	SER	CYS	engineered mutation	UNP Q16881
B	498	CYS	U	SEE REMARK 999	UNP Q16881

- Molecule 2 is a protein called Thioredoxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	105	Total	C	N	O	S	0	0	0
			819	523	128	163	5			
2	D	105	Total	C	N	O	S	0	0	0
			819	523	128	163	5			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	MET	-	expression tag	UNP P10599
C	-9	ARG	-	expression tag	UNP P10599
C	-8	GLY	-	expression tag	UNP P10599
C	-7	SER	-	expression tag	UNP P10599
C	-6	HIS	-	expression tag	UNP P10599
C	-5	HIS	-	expression tag	UNP P10599
C	-4	HIS	-	expression tag	UNP P10599
C	-3	HIS	-	expression tag	UNP P10599
C	-2	HIS	-	expression tag	UNP P10599
C	-1	HIS	-	expression tag	UNP P10599
C	0	GLY	-	expression tag	UNP P10599

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1	SER	-	expression tag	UNP P10599
C	35	SER	CYS	engineered mutation	UNP P10599
C	73	SER	CYS	engineered mutation	UNP P10599
D	-10	MET	-	expression tag	UNP P10599
D	-9	ARG	-	expression tag	UNP P10599
D	-8	GLY	-	expression tag	UNP P10599
D	-7	SER	-	expression tag	UNP P10599
D	-6	HIS	-	expression tag	UNP P10599
D	-5	HIS	-	expression tag	UNP P10599
D	-4	HIS	-	expression tag	UNP P10599
D	-3	HIS	-	expression tag	UNP P10599
D	-2	HIS	-	expression tag	UNP P10599
D	-1	HIS	-	expression tag	UNP P10599
D	0	GLY	-	expression tag	UNP P10599
D	1	SER	-	expression tag	UNP P10599
D	35	SER	CYS	engineered mutation	UNP P10599
D	73	SER	CYS	engineered mutation	UNP P10599

- # FAD

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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		

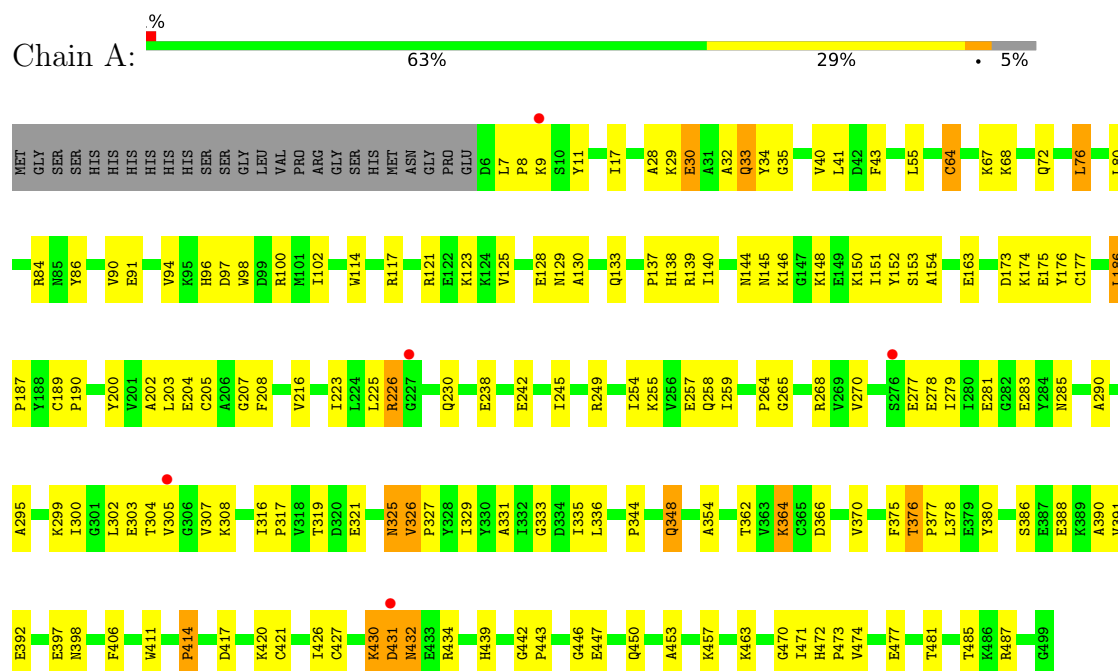
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	190	Total	O	0	0
			190	190		
5	B	202	Total	O	0	0
			202	202		
5	C	17	Total	O	0	0
			17	17		
5	D	16	Total	O	0	0
			16	16		

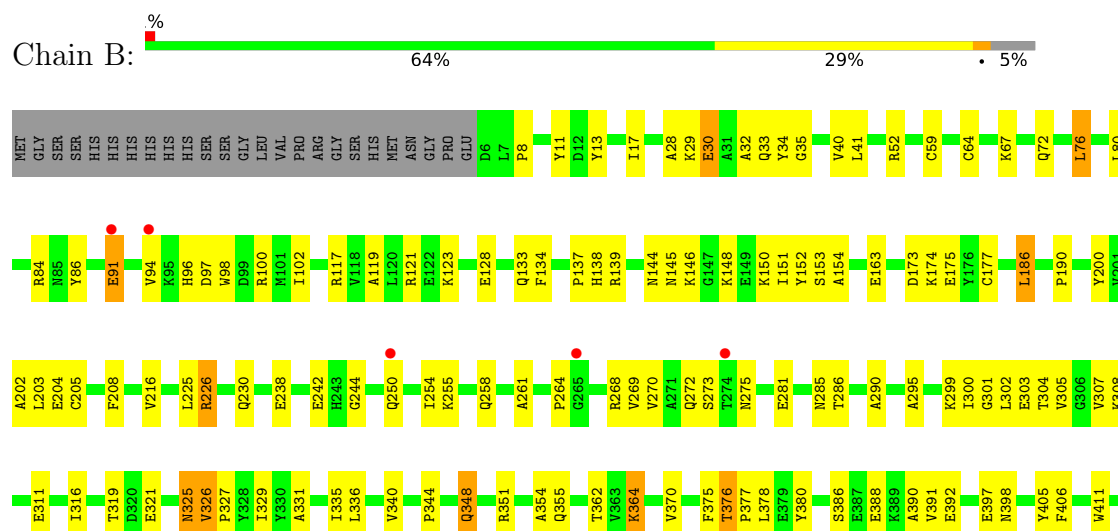
3 Residue-property plots

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: Thioredoxin reductase 1, cytoplasmic

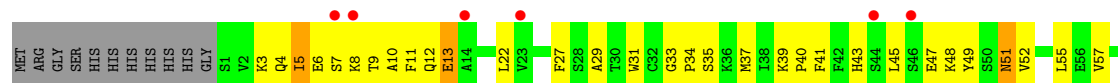


- Molecule 1: Thioredoxin reductase 1, cytoplasmic

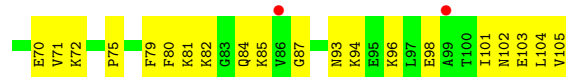
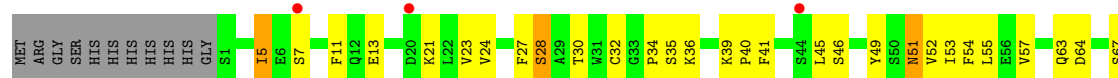




• Molecule 2: Thioredoxin



• Molecule 2: Thioredoxin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	97.25Å 105.03Å 120.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.91 – 2.20 24.91 – 2.20	Depositor EDS
% Data completeness (in resolution range)	84.4 (24.91-2.20) 84.3 (24.91-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.10Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.237 , 0.283 0.237 , 0.283	Depositor DCC
R_{free} test set	11816 reflections (5.96%)	wwPDB-VP
Wilson B-factor (Å ²)	29.3	Xtriage
Anisotropy	0.472	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9821	wwPDB-VP
Average B, all atoms (Å ²)	40.0	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 45.10 % 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 1.3871e-04. 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: FAD, GOL

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	0.41	0/3883	0.98	15/5256 (0.3%)
1	B	0.41	0/3883	0.97	12/5256 (0.2%)
2	C	0.38	0/834	0.79	0/1120
2	D	0.41	0/834	0.84	1/1120 (0.1%)
All	All	0.41	0/9434	0.95	28/12752 (0.2%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	186	LEU	CA-C-N	9.26	128.90	119.82
1	A	186	LEU	C-N-CA	9.26	128.90	119.82
1	B	186	LEU	CA-C-N	9.06	128.70	119.82
1	B	186	LEU	C-N-CA	9.06	128.70	119.82
1	A	326	VAL	CA-C-N	8.39	130.33	119.84
1	A	326	VAL	C-N-CA	8.39	130.33	119.84
1	B	326	VAL	CA-C-N	8.30	130.21	119.84
1	B	326	VAL	C-N-CA	8.30	130.21	119.84
1	B	417	ASP	N-CA-C	6.61	120.37	111.24
2	D	28	SER	N-CA-C	6.58	119.41	109.41
1	B	449	THR	N-CA-C	6.39	118.33	111.36
1	A	417	ASP	N-CA-C	6.35	120.00	111.24
1	A	386	SER	N-CA-C	-6.31	102.39	110.53
1	A	432	ASN	N-CA-C	6.07	118.20	110.61
1	A	414	PRO	N-CA-C	-6.03	106.61	114.03
1	B	386	SER	N-CA-C	-6.01	102.77	110.53
1	B	414	PRO	N-CA-C	-5.71	107.01	114.03
1	A	216	VAL	N-CA-C	5.51	116.41	108.48
1	B	216	VAL	N-CA-C	5.43	116.30	108.48
1	B	230	GLN	N-CA-C	5.39	118.08	111.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	230	GLN	N-CA-C	5.21	117.87	111.82
1	A	189	CYS	CA-C-N	5.18	125.48	119.47
1	A	189	CYS	C-N-CA	5.18	125.48	119.47
1	A	453	ALA	N-CA-C	-5.16	105.73	111.36
1	B	244	GLY	N-CA-C	5.13	121.18	115.08
1	A	9	LYS	N-CA-C	-5.12	106.98	113.18
1	B	453	ALA	N-CA-C	-5.11	105.79	111.36
1	A	226	ARG	N-CA-C	-5.10	103.05	110.24

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	3808	0	3814	170	0
1	B	3808	0	3814	147	0
2	C	819	0	806	49	0
2	D	819	0	806	41	0
3	A	53	0	31	1	0
3	B	53	0	31	0	0
4	A	24	0	16	1	0
4	B	6	0	4	1	0
4	D	6	0	4	0	0
5	A	190	0	0	9	0
5	B	202	0	0	12	0
5	C	17	0	0	1	0
5	D	16	0	0	1	0
All	All	9821	0	9326	368	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 20.

All (368) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:ARG:HH12	1:A:90:VAL:HB	1.20	1.05
1:A:398:ASN:HD22	1:A:430:LYS:HG3	1.26	0.98
1:A:80:LEU:HD23	1:B:80:LEU:HD23	1.50	0.93
2:C:22:LEU:HD11	2:C:79:PHE:HB3	1.58	0.84
1:B:254:ILE:HG13	1:B:255:LYS:HG2	1.59	0.83
1:A:307:VAL:HA	1:A:325:ASN:HD21	1.44	0.83
2:D:23:VAL:HG22	2:D:53:ILE:HB	1.61	0.82
1:B:325:ASN:HD22	1:B:325:ASN:H	1.26	0.82
1:A:398:ASN:ND2	1:A:430:LYS:HG3	1.94	0.81
1:A:84:ARG:NH1	1:A:90:VAL:HB	1.96	0.80
1:B:91:GLU:HG3	1:B:94:VAL:HG22	1.63	0.80
1:B:307:VAL:HA	1:B:325:ASN:HD21	1.45	0.80
1:A:121:ARG:HD3	2:C:67:SER:HB2	1.65	0.79
1:A:325:ASN:HD22	1:A:325:ASN:H	1.28	0.78
2:C:41:PHE:O	2:C:45:LEU:HD13	1.85	0.77
1:B:145:ASN:OD1	1:B:146:LYS:HG3	1.87	0.74
1:A:471:ILE:H	1:B:450:GLN:HE22	1.33	0.74
1:A:145:ASN:OD1	1:A:146:LYS:HG3	1.88	0.73
1:B:311:GLU:HG3	5:B:623:HOH:O	1.88	0.73
1:B:380:TYR:OH	1:B:439:HIS:HD2	1.74	0.71
2:C:81:LYS:HB3	2:C:86:VAL:HG11	1.73	0.71
1:A:471:ILE:H	1:B:450:GLN:NE2	1.89	0.70
1:B:250:GLN:CD	1:B:275:ASN:HD21	1.99	0.70
1:B:250:GLN:NE2	1:B:275:ASN:HD21	1.90	0.69
1:A:254:ILE:HD11	1:A:270:VAL:HG12	1.73	0.69
1:A:380:TYR:OH	1:A:439:HIS:HD2	1.76	0.69
2:C:3:LYS:HB3	2:C:55:LEU:HD23	1.74	0.69
2:C:35:SER:HA	2:C:75:PRO:HG3	1.75	0.69
1:A:325:ASN:HD22	1:A:325:ASN:N	1.91	0.68
1:A:388:GLU:O	1:A:391:VAL:HG22	1.94	0.68
1:A:84:ARG:HH11	1:A:84:ARG:HG3	1.60	0.67
1:B:325:ASN:HD22	1:B:325:ASN:N	1.92	0.67
1:B:434:ARG:HH21	4:B:804:GOL:H11	1.57	0.67
1:A:430:LYS:O	1:A:431:ASP:HB3	1.95	0.67
1:B:117:ARG:O	1:B:121:ARG:HG3	1.94	0.67
1:B:117:ARG:NH2	2:D:63:GLN:HG2	2.10	0.66
1:B:388:GLU:O	1:B:391:VAL:HG22	1.96	0.66
2:D:35:SER:HA	2:D:75:PRO:HG3	1.78	0.66
2:C:84:GLN:O	2:C:86:VAL:HG13	1.96	0.66
1:A:450:GLN:HE22	1:B:471:ILE:H	1.43	0.66
1:A:344:PRO:HG3	1:B:472:HIS:HB2	1.77	0.65
2:D:21:LYS:NZ	2:D:21:LYS:HB3	2.10	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:471:ILE:N	1:B:450:GLN:HE22	1.94	0.65
2:C:4:GLN:O	2:C:6:GLU:HG3	1.97	0.65
1:A:470:GLY:HA2	1:B:450:GLN:HE22	1.60	0.65
1:A:238:GLU:O	1:A:242:GLU:HG3	1.97	0.64
2:D:51:ASN:HD22	2:D:51:ASN:H	1.45	0.64
1:A:268:ARG:HH11	1:A:268:ARG:HB2	1.63	0.63
1:A:463:LYS:HG2	5:A:674:HOH:O	1.97	0.63
1:A:474:VAL:HG13	1:B:447:GLU:CD	2.23	0.63
2:C:51:ASN:HD22	2:C:51:ASN:N	1.96	0.63
2:D:49:TYR:HB3	2:D:51:ASN:HD21	1.64	0.63
1:A:121:ARG:HB3	1:A:121:ARG:HH11	1.64	0.62
1:A:450:GLN:HE22	1:B:470:GLY:HA2	1.64	0.62
2:C:49:TYR:HD1	2:C:101:ILE:HG21	1.65	0.62
2:D:51:ASN:HD22	2:D:51:ASN:N	1.96	0.62
1:B:98:TRP:CZ2	1:B:102:ILE:HD11	2.35	0.61
1:A:133:GLN:C	1:A:140:ILE:HG13	2.25	0.61
2:C:51:ASN:ND2	2:C:52:VAL:HG23	2.16	0.61
1:A:117:ARG:HH21	2:C:63:GLN:HA	1.65	0.61
1:A:268:ARG:HD2	1:A:281:GLU:OE2	2.01	0.61
1:A:457:LYS:HG3	1:B:468:THR:OG1	2.01	0.61
2:D:81:LYS:NZ	2:D:82:LYS:HE3	2.15	0.61
1:A:431:ASP:HB3	4:A:804:GOL:H11	1.82	0.60
2:D:94:LYS:HE3	5:D:344:HOH:O	1.99	0.60
1:B:121:ARG:HG2	5:B:572:HOH:O	2.01	0.60
2:D:101:ILE:O	2:D:105:VAL:HG23	2.00	0.60
1:B:255:LYS:HG3	1:B:270:VAL:HB	1.82	0.60
1:B:426:ILE:HD12	1:B:426:ILE:N	2.17	0.60
1:A:33:GLN:HE21	1:A:33:GLN:HA	1.67	0.59
2:D:21:LYS:HB3	2:D:21:LYS:HZ2	1.68	0.59
2:D:51:ASN:H	2:D:51:ASN:ND2	1.99	0.59
1:B:8:PRO:HG3	1:B:139:ARG:NH1	2.18	0.59
1:A:80:LEU:HD23	1:B:80:LEU:CD2	2.29	0.59
1:B:33:GLN:HE21	1:B:33:GLN:HA	1.67	0.59
1:A:303:GLU:HG2	1:A:304:THR:N	2.18	0.59
1:A:348:GLN:HE21	1:A:348:GLN:HA	1.67	0.58
1:A:7:LEU:HD23	1:A:8:PRO:O	2.04	0.58
1:A:84:ARG:HH12	1:A:90:VAL:CB	2.05	0.58
1:A:96:HIS:HE1	1:B:86:TYR:O	1.85	0.57
2:C:81:LYS:HE3	2:C:82:LYS:HE3	1.85	0.57
1:A:139:ARG:HG2	1:A:139:ARG:HH11	1.68	0.57
1:A:426:ILE:HD12	1:A:426:ILE:N	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:ILE:HG12	1:A:270:VAL:O	2.05	0.57
1:B:305:VAL:HG13	1:B:307:VAL:HG23	1.86	0.57
2:D:34:PRO:HB2	2:D:75:PRO:HD3	1.87	0.57
1:B:139:ARG:HG2	1:B:139:ARG:HH11	1.68	0.57
1:B:348:GLN:HE21	1:B:348:GLN:HA	1.69	0.56
1:A:470:GLY:CA	1:B:450:GLN:HE22	2.18	0.56
1:A:205:CYS:HA	1:A:208:PHE:CE2	2.40	0.56
1:B:119:ALA:O	1:B:123:LYS:HG2	2.04	0.56
1:A:305:VAL:HG13	1:A:307:VAL:HG23	1.87	0.56
1:A:450:GLN:NE2	1:B:471:ILE:H	2.03	0.56
1:B:72:GLN:HG3	1:B:76:LEU:HD22	1.88	0.56
1:A:121:ARG:HH11	1:A:121:ARG:CB	2.19	0.55
1:A:447:GLU:CD	1:B:474:VAL:HG13	2.32	0.55
1:B:250:GLN:CD	1:B:275:ASN:ND2	2.64	0.55
1:B:340:VAL:HG21	1:B:370:VAL:HG21	1.87	0.55
2:C:39:LYS:HB3	2:C:40:PRO:HD3	1.88	0.55
1:B:268:ARG:HD2	1:B:281:GLU:OE2	2.07	0.55
1:B:303:GLU:HG2	1:B:304:THR:N	2.20	0.55
2:C:22:LEU:O	2:C:52:VAL:HG13	2.07	0.55
1:A:446:GLY:HA3	1:B:474:VAL:HG21	1.89	0.55
1:B:11:TYR:HA	1:B:153:SER:OG	2.07	0.55
1:B:411:TRP:CZ2	1:B:443:PRO:HG3	2.42	0.54
2:D:51:ASN:ND2	2:D:52:VAL:HG23	2.23	0.54
1:A:114:TRP:CE2	2:C:74:MET:HE1	2.42	0.54
1:A:268:ARG:HH11	1:A:268:ARG:CB	2.20	0.54
2:D:27:PHE:CD1	2:D:71:VAL:HG21	2.43	0.54
1:A:97:ASP:CG	1:A:100:ARG:HG3	2.32	0.54
1:B:258:GLN:NE2	1:B:261:ALA:HA	2.23	0.54
1:A:470:GLY:HA2	1:B:450:GLN:NE2	2.23	0.54
1:B:308:LYS:HG2	1:B:325:ASN:OD1	2.08	0.54
1:A:255:LYS:HG2	1:A:270:VAL:HB	1.88	0.53
1:A:316:ILE:HD12	1:A:335:ILE:HD12	1.90	0.53
1:A:72:GLN:HG3	1:A:76:LEU:HD22	1.91	0.53
1:A:80:LEU:CD2	1:B:80:LEU:HD23	2.31	0.53
1:B:41:LEU:HD23	1:B:128:GLU:HB3	1.91	0.53
2:C:52:VAL:HG21	2:C:105:VAL:HG21	1.91	0.53
1:A:114:TRP:NE1	2:C:74:MET:HE1	2.25	0.52
1:A:133:GLN:HG3	1:A:300:ILE:C	2.34	0.52
2:D:39:LYS:HB3	2:D:40:PRO:HD3	1.91	0.52
1:A:450:GLN:HE22	1:B:470:GLY:CA	2.23	0.52
1:B:33:GLN:HA	1:B:33:GLN:NE2	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:GLN:HA	1:A:33:GLN:NE2	2.23	0.52
1:A:430:LYS:O	1:A:431:ASP:CB	2.56	0.52
1:A:308:LYS:HG2	1:A:325:ASN:OD1	2.09	0.52
1:B:255:LYS:CG	1:B:270:VAL:HB	2.40	0.52
1:B:316:ILE:HD12	1:B:335:ILE:HD12	1.91	0.52
2:C:49:TYR:CD1	2:C:101:ILE:HG21	2.45	0.51
1:B:238:GLU:O	1:B:242:GLU:HG3	2.10	0.51
2:D:49:TYR:HB3	2:D:51:ASN:ND2	2.25	0.51
2:C:22:LEU:HD22	2:C:104:LEU:HB2	1.93	0.51
1:B:205:CYS:HA	1:B:208:PHE:CE2	2.44	0.51
1:A:17:ILE:HD12	1:A:28:ALA:HB2	1.93	0.51
2:C:11:PHE:CD2	2:C:65:VAL:HG13	2.46	0.51
1:B:17:ILE:HD12	1:B:28:ALA:HB2	1.92	0.50
1:B:325:ASN:H	1:B:325:ASN:ND2	2.03	0.50
2:D:41:PHE:O	2:D:45:LEU:HB2	2.10	0.50
1:A:370:VAL:HG13	1:B:469:ILE:HB	1.92	0.50
1:B:226:ARG:NH1	1:B:226:ARG:HG3	2.26	0.50
2:C:7:SER:HB3	2:C:10:ALA:CB	2.42	0.50
1:A:117:ARG:HG3	1:A:117:ARG:HH11	1.77	0.50
1:A:145:ASN:HB2	5:A:677:HOH:O	2.12	0.50
1:A:472:HIS:HB2	1:B:344:PRO:HG3	1.92	0.50
1:B:226:ARG:HH11	1:B:226:ARG:CG	2.25	0.50
1:A:98:TRP:CZ2	1:A:102:ILE:HD11	2.47	0.50
1:A:123:LYS:HB2	1:A:125:VAL:HG23	1.93	0.50
1:A:7:LEU:HD23	1:A:7:LEU:C	2.36	0.50
1:A:474:VAL:O	1:A:477:GLU:HG2	2.12	0.50
1:B:474:VAL:O	1:B:477:GLU:HG2	2.12	0.50
2:C:81:LYS:HB2	2:C:86:VAL:HG21	1.92	0.50
2:D:84:GLN:NE2	2:D:85:LYS:H	2.09	0.50
1:B:91:GLU:HG3	1:B:94:VAL:CG2	2.38	0.50
1:B:264:PRO:HB2	1:B:285:ASN:HB3	1.94	0.49
2:D:5:ILE:CG1	2:D:57:VAL:HG22	2.41	0.49
1:B:84:ARG:HB3	1:B:84:ARG:NH1	2.28	0.49
1:A:67:LYS:C	1:A:67:LYS:HD2	2.37	0.49
1:A:255:LYS:CG	1:A:270:VAL:HB	2.42	0.49
1:A:257:GLU:OE1	1:A:268:ARG:NH2	2.45	0.49
1:A:406:PHE:CZ	1:A:421:CYS:HB3	2.47	0.49
1:A:411:TRP:C	1:A:414:PRO:HD2	2.37	0.49
1:B:97:ASP:CG	1:B:100:ARG:HG3	2.37	0.49
1:B:137:PRO:O	1:B:139:ARG:HG3	2.13	0.49
2:C:8:LYS:HB3	2:C:64:ASP:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ASN:ND2	1:A:148:LYS:HB3	2.27	0.49
2:D:5:ILE:HD12	2:D:5:ILE:C	2.37	0.49
1:A:8:PRO:HG2	1:A:153:SER:CB	2.41	0.49
1:A:121:ARG:CB	1:A:121:ARG:NH1	2.76	0.49
1:A:283:GLU:N	5:A:686:HOH:O	2.40	0.48
1:A:207:GLY:HA2	1:A:245:ILE:HD11	1.95	0.48
1:A:254:ILE:HD11	1:A:270:VAL:CG1	2.41	0.48
1:B:406:PHE:CZ	1:B:421:CYS:HB3	2.48	0.48
1:A:150:LYS:HD3	1:A:152:TYR:OH	2.14	0.48
1:A:397:GLU:CD	1:A:397:GLU:H	2.22	0.48
1:A:30:GLU:O	1:A:33:GLN:HB3	2.14	0.48
1:A:137:PRO:O	1:A:139:ARG:HG3	2.13	0.48
1:A:450:GLN:HE22	1:B:471:ILE:N	2.09	0.48
1:B:258:GLN:HE21	1:B:261:ALA:HA	1.78	0.48
1:B:299:LYS:NZ	1:B:299:LYS:HB3	2.29	0.48
1:B:493:LEU:CD1	1:B:493:LEU:N	2.76	0.48
1:A:203:LEU:HD12	1:A:225:LEU:HD21	1.96	0.48
1:B:203:LEU:HD12	1:B:225:LEU:HD21	1.94	0.48
1:A:30:GLU:HB3	1:A:354:ALA:HB3	1.96	0.48
1:A:117:ARG:NE	2:C:63:GLN:HB3	2.29	0.48
1:B:286:THR:HG23	5:B:543:HOH:O	2.14	0.48
1:B:13:TYR:O	1:B:154:ALA:HA	2.13	0.48
1:A:268:ARG:CB	1:A:268:ARG:NH1	2.77	0.47
1:A:11:TYR:HE1	5:A:661:HOH:O	1.97	0.47
2:D:21:LYS:NZ	2:D:51:ASN:O	2.46	0.47
1:A:114:TRP:CD1	2:C:74:MET:HE1	2.49	0.47
1:B:67:LYS:HD2	1:B:67:LYS:C	2.39	0.47
1:A:139:ARG:HG2	1:A:139:ARG:NH1	2.29	0.47
1:A:299:LYS:NZ	1:A:299:LYS:HB3	2.28	0.47
1:A:431:ASP:O	1:A:432:ASN:OD1	2.32	0.47
1:B:30:GLU:OE1	1:B:30:GLU:HA	2.14	0.47
1:B:272:GLN:HG2	1:B:273:SER:N	2.28	0.47
1:B:121:ARG:HH11	1:B:121:ARG:HB3	1.80	0.47
2:C:69:CYS:HB3	2:C:78:GLN:HE22	1.80	0.47
1:A:450:GLN:NE2	1:B:477:GLU:HA	2.29	0.47
1:B:59:CYS:HB3	5:B:653:HOH:O	2.15	0.47
1:B:186:LEU:CD2	1:B:190:PRO:HG3	2.45	0.47
1:B:397:GLU:CD	1:B:397:GLU:H	2.22	0.47
1:B:305:VAL:HG21	1:B:329:ILE:HD11	1.96	0.47
2:C:7:SER:HB3	2:C:10:ALA:HB2	1.96	0.47
2:D:81:LYS:HZ2	2:D:82:LYS:HE3	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:ASP:O	1:A:177:CYS:HB2	2.15	0.47
1:A:30:GLU:OE1	1:A:30:GLU:HA	2.14	0.46
1:B:30:GLU:O	1:B:33:GLN:HB3	2.15	0.46
2:C:12:GLN:NE2	2:C:12:GLN:HA	2.30	0.46
1:A:86:TYR:O	1:B:96:HIS:HE1	1.98	0.46
1:B:144:ASN:ND2	1:B:148:LYS:HB3	2.30	0.46
1:B:173:ASP:O	1:B:177:CYS:HB2	2.15	0.46
2:C:43:HIS:O	2:C:47:GLU:HG2	2.15	0.46
1:A:138:HIS:O	1:A:153:SER:HA	2.16	0.46
1:A:305:VAL:HG21	1:A:329:ILE:HD11	1.97	0.46
1:B:29:LYS:NZ	5:B:624:HOH:O	2.44	0.46
1:B:30:GLU:HB3	1:B:354:ALA:HB3	1.97	0.46
1:B:138:HIS:O	1:B:153:SER:HA	2.15	0.46
2:C:72:LYS:HG2	5:C:114:HOH:O	2.15	0.46
1:A:32:ALA:O	1:A:35:GLY:N	2.48	0.46
1:A:64:CYS:SG	1:B:472:HIS:HD2	2.38	0.46
1:A:121:ARG:HE	2:C:67:SER:HA	1.80	0.46
1:A:362:THR:O	1:A:364:LYS:HD2	2.15	0.46
1:B:139:ARG:NH1	1:B:139:ARG:HG2	2.29	0.46
1:B:133:GLN:NE2	5:B:530:HOH:O	2.35	0.46
2:C:29:ALA:HB1	2:C:31:TRP:CE2	2.50	0.46
2:C:81:LYS:HB3	2:C:86:VAL:CG1	2.44	0.46
1:A:176:TYR:CE1	1:A:258:GLN:HB2	2.51	0.45
1:A:326:VAL:HA	1:A:327:PRO:HD3	1.82	0.45
1:B:150:LYS:HD3	1:B:152:TYR:OH	2.16	0.45
2:C:4:GLN:HA	2:C:4:GLN:NE2	2.31	0.45
1:A:117:ARG:HG3	1:A:117:ARG:NH1	2.32	0.45
1:A:173:ASP:OD1	1:A:174:LYS:N	2.49	0.45
1:A:186:LEU:CD2	1:A:190:PRO:HG3	2.46	0.45
1:A:487:ARG:HG3	1:A:487:ARG:HH11	1.81	0.45
2:D:79:PHE:HB2	2:D:87:GLY:HA3	1.99	0.45
1:A:485:THR:HB	5:A:508:HOH:O	2.16	0.45
1:B:496:GLY:H	1:B:499:GLY:C	2.24	0.45
2:C:22:LEU:HD23	2:C:105:VAL:CG2	2.47	0.45
1:B:391:VAL:HG23	1:B:392:GLU:N	2.31	0.45
1:A:319:THR:C	1:A:321:GLU:H	2.25	0.45
1:A:411:TRP:CZ2	1:A:443:PRO:HG3	2.52	0.45
1:A:55:LEU:HD23	1:A:117:ARG:HG2	1.97	0.45
1:A:117:ARG:NH2	2:C:63:GLN:HG2	2.30	0.45
1:B:173:ASP:OD1	1:B:174:LYS:N	2.50	0.45
1:A:41:LEU:HD23	1:A:128:GLU:HB3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:ARG:NH1	1:A:84:ARG:HG3	2.29	0.45
1:A:151:ILE:N	1:A:151:ILE:HD12	2.32	0.44
1:A:249:ARG:HB3	5:A:675:HOH:O	2.16	0.44
1:B:32:ALA:O	1:B:35:GLY:N	2.49	0.44
1:B:139:ARG:NH2	1:B:151:ILE:HG21	2.32	0.44
1:B:151:ILE:HG22	5:B:552:HOH:O	2.17	0.44
1:A:11:TYR:HA	1:A:153:SER:OG	2.18	0.44
1:A:139:ARG:NH2	1:A:151:ILE:HG21	2.33	0.44
1:B:362:THR:O	1:B:364:LYS:HD2	2.16	0.44
2:C:64:ASP:O	2:C:65:VAL:C	2.60	0.44
1:A:163:GLU:HG2	1:A:295:ALA:HA	1.99	0.44
2:D:24:VAL:HG22	2:D:79:PHE:CD2	2.53	0.44
1:A:265:GLY:O	1:A:285:ASN:HA	2.17	0.44
1:A:390:ALA:CB	1:A:426:ILE:HG21	2.48	0.44
1:A:450:GLN:NE2	1:B:470:GLY:HA2	2.32	0.44
1:A:450:GLN:HE21	1:B:477:GLU:HA	1.82	0.44
1:B:269:VAL:O	1:B:281:GLU:HA	2.16	0.44
1:B:398:ASN:HD22	1:B:430:LYS:HG3	1.82	0.44
1:B:487:ARG:HG3	1:B:487:ARG:HH11	1.82	0.44
2:C:5:ILE:HD12	2:C:5:ILE:C	2.43	0.44
2:C:97:LEU:O	2:C:101:ILE:HG13	2.18	0.44
2:D:30:THR:O	2:D:36:LYS:HD3	2.18	0.44
1:A:391:VAL:HG23	1:A:392:GLU:N	2.31	0.44
1:B:331:ALA:C	1:B:336:LEU:HD21	2.42	0.44
2:C:22:LEU:HD23	2:C:105:VAL:HG23	1.99	0.44
1:B:376:THR:O	1:B:377:PRO:C	2.60	0.44
1:B:151:ILE:HD12	1:B:151:ILE:N	2.32	0.43
1:B:226:ARG:NH1	1:B:226:ARG:CG	2.80	0.43
1:B:325:ASN:N	1:B:325:ASN:ND2	2.63	0.43
1:A:225:LEU:O	1:A:226:ARG:C	2.61	0.43
1:A:366:ASP:HB3	5:A:541:HOH:O	2.17	0.43
1:B:431:ASP:OD1	1:B:432:ASN:ND2	2.52	0.43
1:A:11:TYR:HB3	1:A:154:ALA:HA	2.01	0.43
1:A:43:PHE:HA	5:A:514:HOH:O	2.18	0.43
2:D:93:ASN:OD1	2:D:96:LYS:N	2.41	0.43
1:A:32:ALA:C	1:A:34:TYR:N	2.76	0.43
1:A:40:VAL:O	1:A:40:VAL:HG23	2.18	0.43
1:A:29:LYS:HD2	1:A:123:LYS:NZ	2.33	0.43
1:B:117:ARG:HB3	5:B:664:HOH:O	2.18	0.43
1:B:405:TYR:CD1	1:B:492:ILE:HG13	2.53	0.43
1:B:498:CYS:SG	2:C:34:PRO:HD2	2.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:5:ILE:HG13	2:D:57:VAL:HG22	2.01	0.43
1:B:40:VAL:HG23	1:B:40:VAL:O	2.18	0.43
2:C:48:LYS:HE2	2:C:49:TYR:HE2	1.82	0.43
1:B:378:LEU:HD12	1:B:378:LEU:HA	1.84	0.43
1:B:390:ALA:CB	1:B:426:ILE:HG21	2.48	0.43
2:D:5:ILE:HG21	2:D:55:LEU:HD13	2.00	0.43
2:D:103:GLU:O	2:D:103:GLU:HG2	2.17	0.43
1:B:134:PHE:HB2	1:B:301:GLY:O	2.18	0.43
1:A:277:GLU:O	1:A:278:GLU:C	2.62	0.42
1:B:319:THR:C	1:B:321:GLU:H	2.27	0.42
1:B:326:VAL:HA	1:B:327:PRO:HD3	1.82	0.42
2:C:27:PHE:CD1	2:C:57:VAL:HB	2.53	0.42
1:A:277:GLU:O	1:A:279:ILE:HG13	2.19	0.42
1:B:427:CYS:HA	1:B:434:ARG:O	2.20	0.42
1:B:425:ILE:C	1:B:426:ILE:HD12	2.44	0.42
1:A:90:VAL:HG12	1:A:91:GLU:N	2.35	0.42
1:A:378:LEU:HD11	1:A:442:GLY:HA2	2.00	0.42
1:A:420:LYS:HA	1:A:420:LYS:HD3	1.88	0.42
2:C:4:GLN:HA	2:C:4:GLN:HE21	1.85	0.42
2:C:9:THR:O	2:C:13:GLU:HB2	2.20	0.42
1:B:202:ALA:HB2	1:B:290:ALA:HB3	2.01	0.42
1:A:202:ALA:HB2	1:A:290:ALA:HB3	2.01	0.42
1:A:398:ASN:ND2	1:A:430:LYS:CG	2.75	0.42
1:B:200:TYR:O	1:B:204:GLU:HG3	2.20	0.42
1:B:378:LEU:HD11	1:B:442:GLY:HA2	2.00	0.42
1:A:223:ILE:HD11	1:A:226:ARG:HG3	2.00	0.41
1:A:333:GLY:HA3	3:A:600:FAD:O2P	2.19	0.41
1:A:427:CYS:HA	1:A:434:ARG:O	2.20	0.41
1:A:431:ASP:OD2	1:A:432:ASN:ND2	2.52	0.41
2:D:46:SER:HA	2:D:54:PHE:CZ	2.55	0.41
1:A:144:ASN:HD21	1:A:148:LYS:HB3	1.84	0.41
1:A:200:TYR:O	1:A:204:GLU:HG3	2.20	0.41
1:B:91:GLU:CG	1:B:94:VAL:HG22	2.43	0.41
2:D:28:SER:HB2	2:D:35:SER:OG	2.20	0.41
1:A:94:VAL:HA	5:A:628:HOH:O	2.20	0.41
1:A:128:GLU:O	1:A:130:ALA:N	2.52	0.41
1:A:144:ASN:HD21	1:A:148:LYS:CB	2.33	0.41
2:C:33:GLY:O	2:C:37:MET:HG2	2.20	0.41
1:A:331:ALA:C	1:A:336:LEU:HD21	2.46	0.41
1:A:33:GLN:HE21	1:A:33:GLN:CA	2.28	0.41
1:A:226:ARG:HH11	1:A:226:ARG:HB2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:ILE:HD12	1:A:283:GLU:HG3	2.02	0.41
1:A:376:THR:O	1:A:377:PRO:C	2.63	0.41
1:B:163:GLU:HG2	1:B:295:ALA:HA	2.03	0.41
1:B:417:ASP:HB3	1:B:420:LYS:HG3	2.03	0.41
2:D:80:PHE:HA	2:D:84:GLN:O	2.21	0.41
1:A:302:LEU:O	1:A:305:VAL:HG12	2.21	0.41
1:B:186:LEU:HD23	1:B:190:PRO:HG3	2.02	0.41
2:D:104:LEU:O	2:D:105:VAL:C	2.62	0.41
1:B:268:ARG:HH11	1:B:268:ARG:HB2	1.84	0.41
1:A:67:LYS:NZ	1:A:375:PHE:CD1	2.88	0.41
1:A:97:ASP:OD1	1:A:100:ARG:HG3	2.21	0.41
1:B:493:LEU:N	1:B:493:LEU:HD12	2.35	0.41
1:A:8:PRO:HG2	1:A:153:SER:HB2	2.02	0.41
1:A:316:ILE:HA	1:A:317:PRO:HD3	1.90	0.41
1:B:67:LYS:NZ	1:B:375:PHE:CD1	2.89	0.41
1:B:300:ILE:HG13	1:B:302:LEU:HG	2.03	0.41
1:B:391:VAL:CG2	1:B:392:GLU:N	2.84	0.41
1:B:487:ARG:HG3	5:B:689:HOH:O	2.20	0.41
2:C:8:LYS:HB3	2:C:64:ASP:CB	2.51	0.41
2:C:64:ASP:O	2:C:67:SER:OG	2.36	0.41
2:D:98:GLU:O	2:D:102:ASN:ND2	2.54	0.41
1:A:186:LEU:HA	1:A:187:PRO:HD3	1.70	0.41
2:D:5:ILE:HD13	2:D:11:PHE:HB2	2.02	0.41
2:D:24:VAL:HG21	2:D:101:ILE:HG12	2.03	0.41
2:D:81:LYS:HZ1	2:D:82:LYS:HE3	1.81	0.41
1:A:150:LYS:C	1:A:151:ILE:HD12	2.46	0.40
1:A:259:ILE:HG12	1:A:268:ARG:NH1	2.35	0.40
1:A:344:PRO:CG	1:B:472:HIS:HB2	2.49	0.40
1:B:268:ARG:HD2	1:B:281:GLU:CD	2.45	0.40
1:B:355:GLN:NE2	5:B:676:HOH:O	2.53	0.40
2:D:46:SER:HA	2:D:54:PHE:CE1	2.56	0.40
1:A:98:TRP:O	1:A:102:ILE:HG12	2.21	0.40
2:D:32:CYS:SG	2:D:34:PRO:HD2	2.62	0.40
1:A:481:THR:HG21	1:B:351:ARG:HH12	1.86	0.40
1:A:68:LYS:HE2	5:B:558:HOH:O	2.22	0.40
1:B:32:ALA:C	1:B:34:TYR:N	2.77	0.40
1:B:370:VAL:HG23	5:B:668:HOH:O	2.21	0.40
2:D:70:GLU:OE2	2:D:72:LYS:HE3	2.22	0.40
1:B:331:ALA:O	1:B:336:LEU:HD21	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	492/519 (95%)	462 (94%)	25 (5%)	5 (1%)	12	11
1	B	492/519 (95%)	464 (94%)	26 (5%)	2 (0%)	30	34
2	C	103/116 (89%)	90 (87%)	12 (12%)	1 (1%)	12	11
2	D	103/116 (89%)	100 (97%)	3 (3%)	0	100	100
All	All	1190/1270 (94%)	1116 (94%)	66 (6%)	8 (1%)	18	19

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	431	ASP
2	C	65	VAL
1	A	129	ASN
1	B	430	LYS
1	A	33	GLN
1	A	430	LYS
1	B	473	PRO
1	A	473	PRO

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	405/426 (95%)	396 (98%)	9 (2%)	45	61
1	B	405/426 (95%)	394 (97%)	11 (3%)	39	53

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	92/101 (91%)	89 (97%)	3 (3%)	33	45
2	D	92/101 (91%)	86 (94%)	6 (6%)	15	18
All	All	994/1054 (94%)	965 (97%)	29 (3%)	37	51

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLU
1	A	64	CYS
1	A	76	LEU
1	A	175	GLU
1	A	264	PRO
1	A	325	ASN
1	A	348	GLN
1	A	364	LYS
1	A	376	THR
1	B	30	GLU
1	B	52	ARG
1	B	64	CYS
1	B	76	LEU
1	B	91	GLU
1	B	175	GLU
1	B	226	ARG
1	B	325	ASN
1	B	348	GLN
1	B	364	LYS
1	B	376	THR
2	C	5	ILE
2	C	13	GLU
2	C	51	ASN
2	D	5	ILE
2	D	7	SER
2	D	13	GLU
2	D	51	ASN
2	D	64	ASP
2	D	67	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	96	HIS
1	A	106	GLN
1	A	107	ASN
1	A	129	ASN
1	A	138	HIS
1	A	275	ASN
1	A	325	ASN
1	A	348	GLN
1	A	355	GLN
1	A	398	ASN
1	A	428	ASN
1	A	432	ASN
1	A	439	HIS
1	A	450	GLN
1	B	33	GLN
1	B	81	GLN
1	B	96	HIS
1	B	107	ASN
1	B	129	ASN
1	B	138	HIS
1	B	250	GLN
1	B	258	GLN
1	B	275	ASN
1	B	325	ASN
1	B	348	GLN
1	B	355	GLN
1	B	398	ASN
1	B	439	HIS
1	B	450	GLN
1	B	472	HIS
2	C	4	GLN
2	C	12	GLN
2	C	51	ASN
2	C	63	GLN
2	C	78	GLN
2	C	102	ASN
2	D	4	GLN
2	D	12	GLN
2	D	51	ASN
2	D	84	GLN
2	D	102	ASN

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$
4	GOL	D	803	-	5,5,5	4.84	5 (100%)	5,5,5	6.11	3 (60%)
4	GOL	A	802	-	5,5,5	4.77	5 (100%)	5,5,5	6.14	3 (60%)
3	FAD	A	600	-	58,58,58	1.55	9 (15%)	85,89,89	0.81	2 (2%)
4	GOL	A	804	-	5,5,5	4.80	5 (100%)	5,5,5	6.12	3 (60%)
4	GOL	B	804	-	5,5,5	4.82	5 (100%)	5,5,5	6.12	3 (60%)
4	GOL	A	803	-	5,5,5	4.83	5 (100%)	5,5,5	6.13	3 (60%)
3	FAD	B	600	-	58,58,58	1.46	8 (13%)	85,89,89	0.86	2 (2%)
4	GOL	A	801	-	5,5,5	4.74	5 (100%)	5,5,5	6.16	3 (60%)

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
4	GOL	D	803	-	-	2/4/4/4	-
4	GOL	A	802	-	-	2/4/4/4	-
3	FAD	A	600	-	-	5/34/50/50	0/6/6/6
4	GOL	A	804	-	-	2/4/4/4	-
4	GOL	B	804	-	-	3/4/4/4	-
4	GOL	A	803	-	-	2/4/4/4	-
3	FAD	B	600	-	-	5/34/50/50	0/6/6/6
4	GOL	A	801	-	-	2/4/4/4	-

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	803	GOL	C3-C2	-8.11	1.20	1.51
4	D	803	GOL	C3-C2	-8.11	1.20	1.51
4	B	804	GOL	C3-C2	-8.06	1.21	1.51
4	A	804	GOL	C3-C2	-8.05	1.21	1.51
4	A	802	GOL	C3-C2	-8.01	1.21	1.51
4	A	801	GOL	C3-C2	-7.99	1.21	1.51
3	B	600	FAD	C4X-N5	5.39	1.42	1.30
3	A	600	FAD	C4X-N5	5.23	1.42	1.30
4	A	801	GOL	O1-C1	4.67	1.62	1.42
4	A	803	GOL	O1-C1	4.60	1.61	1.42
4	B	804	GOL	O1-C1	4.51	1.61	1.42
4	A	802	GOL	O1-C1	4.48	1.61	1.42
4	D	803	GOL	O1-C1	4.46	1.61	1.42
4	A	804	GOL	O1-C1	4.29	1.60	1.42
3	A	600	FAD	C9-C8	4.17	1.45	1.39
3	B	600	FAD	C4A-N3A	3.76	1.41	1.34
3	B	600	FAD	C9-C8	3.76	1.44	1.39
3	A	600	FAD	C4A-N3A	3.69	1.41	1.34
3	A	600	FAD	C9A-C5X	3.53	1.46	1.41
4	A	804	GOL	O3-C3	3.49	1.57	1.42
4	A	802	GOL	O3-C3	3.46	1.56	1.42
4	D	803	GOL	O3-C3	3.45	1.56	1.42
4	A	803	GOL	O3-C3	3.39	1.56	1.42
4	B	804	GOL	O3-C3	3.38	1.56	1.42
4	A	804	GOL	C1-C2	-3.31	1.39	1.51
4	A	801	GOL	O3-C3	3.29	1.56	1.42
4	B	804	GOL	C1-C2	-3.15	1.39	1.51
4	D	803	GOL	C1-C2	-3.15	1.39	1.51
4	D	803	GOL	O2-C2	-3.13	1.34	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	802	GOL	C1-C2	-3.12	1.39	1.51
4	B	804	GOL	O2-C2	-3.09	1.34	1.43
4	A	803	GOL	C1-C2	-3.08	1.40	1.51
3	A	600	FAD	C10-N1	3.05	1.39	1.33
3	A	600	FAD	C9A-N10	3.02	1.46	1.41
4	A	804	GOL	O2-C2	-2.98	1.34	1.43
4	A	801	GOL	C1-C2	-2.97	1.40	1.51
3	B	600	FAD	C10-N1	2.94	1.39	1.33
4	A	803	GOL	O2-C2	-2.93	1.34	1.43
4	A	802	GOL	O2-C2	-2.81	1.35	1.43
3	A	600	FAD	P-O3P	2.80	1.62	1.59
4	A	801	GOL	O2-C2	-2.67	1.35	1.43
3	B	600	FAD	C9A-C5X	2.64	1.45	1.41
3	B	600	FAD	C6-C5X	2.55	1.43	1.40
3	A	600	FAD	C6-C5X	2.50	1.43	1.40
3	B	600	FAD	P-O3P	2.47	1.62	1.59
3	B	600	FAD	C9A-N10	2.21	1.45	1.41
3	A	600	FAD	C2A-N1A	2.06	1.37	1.33

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	801	GOL	O3-C3-C2	11.24	160.96	110.38
4	A	803	GOL	O3-C3-C2	11.20	160.78	110.38
4	A	802	GOL	O3-C3-C2	11.17	160.69	110.38
4	B	804	GOL	O3-C3-C2	11.16	160.64	110.38
4	A	804	GOL	O3-C3-C2	11.07	160.21	110.38
4	D	803	GOL	O3-C3-C2	11.07	160.19	110.38
4	A	804	GOL	O2-C2-C3	7.31	139.45	109.18
4	D	803	GOL	O2-C2-C3	7.19	138.96	109.18
4	A	802	GOL	O2-C2-C3	7.13	138.71	109.18
4	A	801	GOL	O2-C2-C3	7.11	138.61	109.18
4	B	804	GOL	O2-C2-C3	7.10	138.58	109.18
4	A	803	GOL	O2-C2-C3	7.10	138.57	109.18
4	A	801	GOL	O1-C1-C2	3.49	126.07	110.38
4	A	802	GOL	O1-C1-C2	3.47	126.00	110.38
4	D	803	GOL	O1-C1-C2	3.39	125.66	110.38
4	B	804	GOL	O1-C1-C2	3.38	125.59	110.38
4	A	803	GOL	O1-C1-C2	3.38	125.57	110.38
4	A	804	GOL	O1-C1-C2	3.27	125.08	110.38
3	A	600	FAD	C4X-C10-N10	2.30	119.77	116.48
3	B	600	FAD	C4X-C10-N10	2.26	119.72	116.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	600	FAD	C2B-C1B-N9A	2.22	118.81	113.30
3	A	600	FAD	C5'-C4'-C3'	-2.00	108.44	112.22

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	801	GOL	C1-C2-C3-O3
4	A	802	GOL	C1-C2-C3-O3
4	A	804	GOL	C1-C2-C3-O3
4	A	803	GOL	O1-C1-C2-C3
4	A	803	GOL	C1-C2-C3-O3
4	B	804	GOL	C1-C2-C3-O3
4	D	803	GOL	O1-C1-C2-C3
4	D	803	GOL	C1-C2-C3-O3
4	A	804	GOL	O1-C1-C2-C3
4	B	804	GOL	O1-C1-C2-C3
4	A	801	GOL	O1-C1-C2-O2
4	B	804	GOL	O1-C1-C2-O2
4	A	802	GOL	O1-C1-C2-O2
3	A	600	FAD	O4B-C4B-C5B-O5B
3	A	600	FAD	PA-O3P-P-O5'
3	B	600	FAD	PA-O3P-P-O5'
3	A	600	FAD	P-O3P-PA-O1A
3	B	600	FAD	P-O3P-PA-O1A
3	A	600	FAD	C3B-C4B-C5B-O5B
3	A	600	FAD	P-O3P-PA-O2A
3	B	600	FAD	P-O3P-PA-O2A
3	B	600	FAD	O4B-C4B-C5B-O5B
3	B	600	FAD	C2B-C1B-N9A-C8A

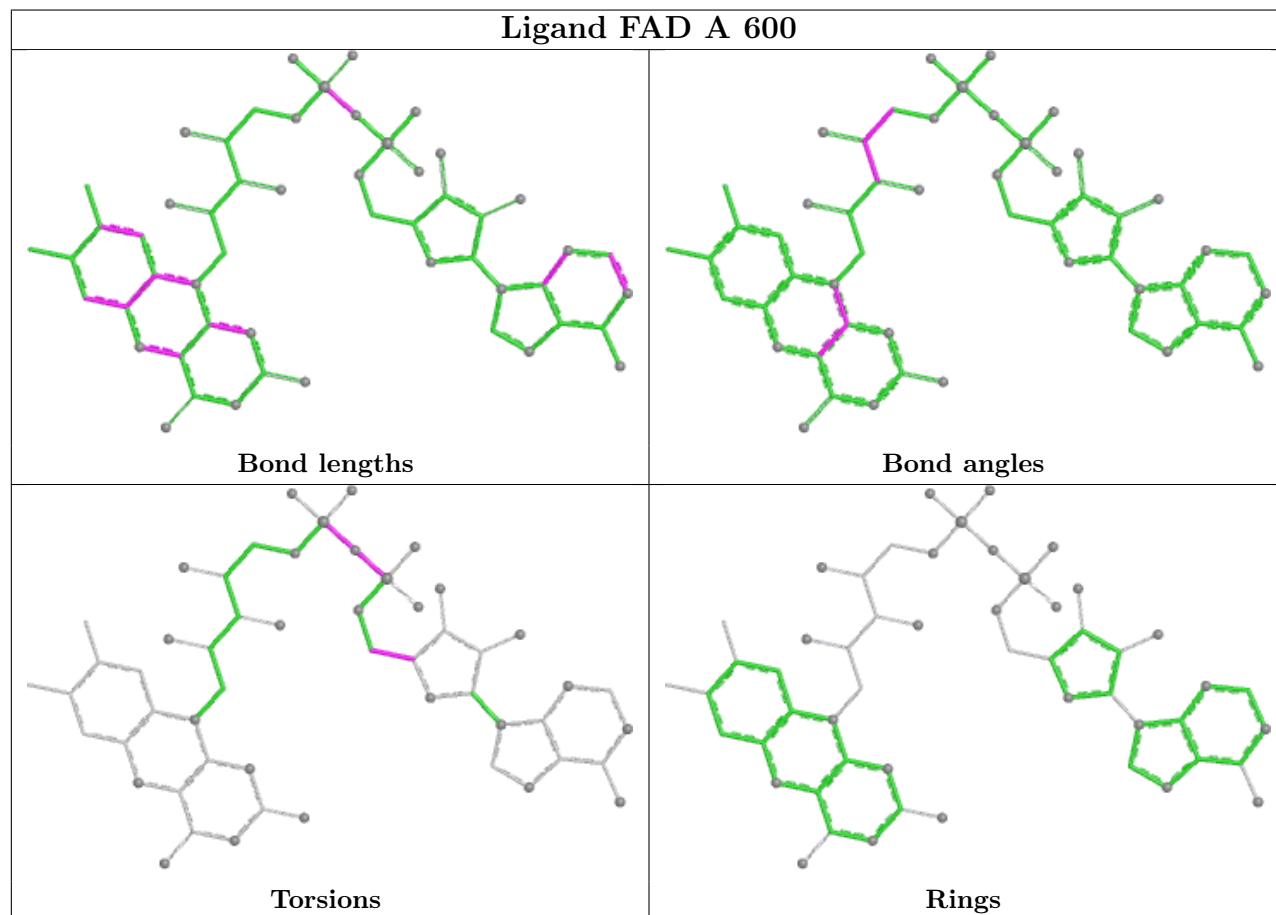
There are no ring outliers.

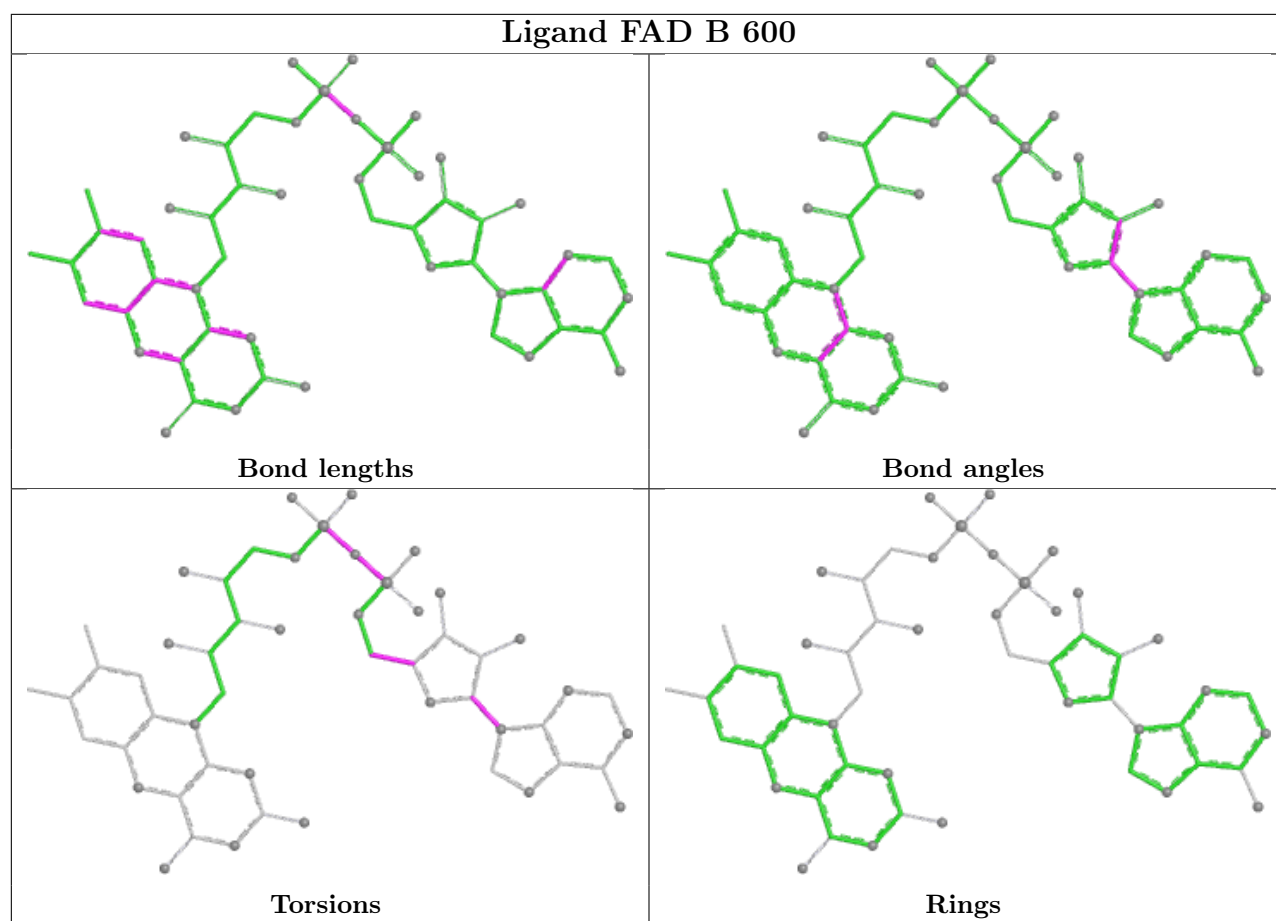
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	600	FAD	1	0
4	A	804	GOL	1	0
4	B	804	GOL	1	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	494/519 (95%)	0.33	5 (1%) 79 77	22, 37, 57, 88	0
1	B	494/519 (95%)	0.27	7 (1%) 73 71	21, 35, 58, 80	0
2	C	105/116 (90%)	1.05	13 (12%) 8 6	30, 54, 68, 74	0
2	D	105/116 (90%)	0.64	5 (4%) 35 32	29, 42, 63, 69	0
All	All	1198/1270 (94%)	0.40	30 (2%) 58 55	21, 38, 63, 88	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	497	SER	4.1
2	C	66	ALA	4.1
2	C	67	SER	3.8
2	C	7	SER	3.6
1	A	305	VAL	3.3
1	A	431	ASP	3.3
2	C	86	VAL	3.1
1	B	250	GLN	3.1
1	B	94	VAL	3.0
2	C	14	ALA	2.9
2	D	44	SER	2.9
1	B	431	ASP	2.9
2	C	83	GLY	2.7
2	C	65	VAL	2.7
1	B	91	GLU	2.5
1	A	9	LYS	2.4
1	B	265	GLY	2.3
1	B	274	THR	2.3
2	C	23	VAL	2.3
1	A	276	SER	2.3
2	C	71	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
2	C	46	SER	2.3
1	A	227	GLY	2.2
2	D	20	ASP	2.2
2	D	7	SER	2.2
2	D	86	VAL	2.2
2	C	8	LYS	2.1
2	C	44	SER	2.1
2	C	88	GLU	2.0
2	D	99	ALA	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](#)

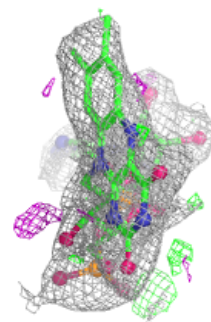
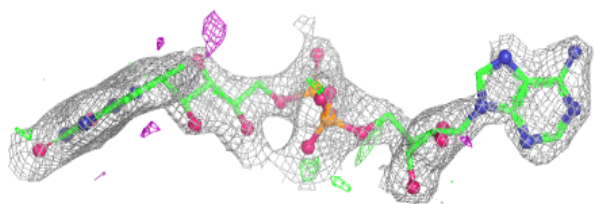
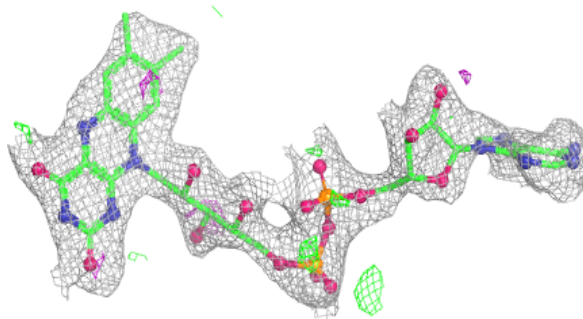
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	804	6/6	0.69	0.17	59,65,66,68	0
4	GOL	A	801	6/6	0.73	0.15	60,61,62,62	0
4	GOL	A	804	6/6	0.78	0.15	52,54,54,55	0
4	GOL	A	802	6/6	0.82	0.15	57,60,61,62	0
4	GOL	D	803	6/6	0.87	0.12	68,69,70,70	0
4	GOL	A	803	6/6	0.89	0.08	55,55,56,58	0
3	FAD	A	600	53/53	0.92	0.10	32,39,42,42	0
3	FAD	B	600	53/53	0.96	0.07	23,29,32,34	0

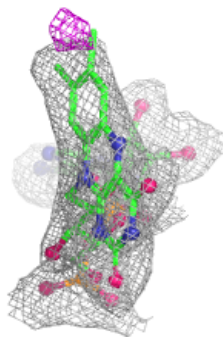
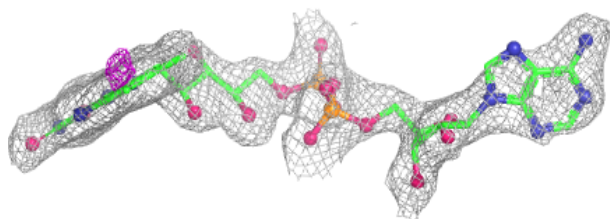
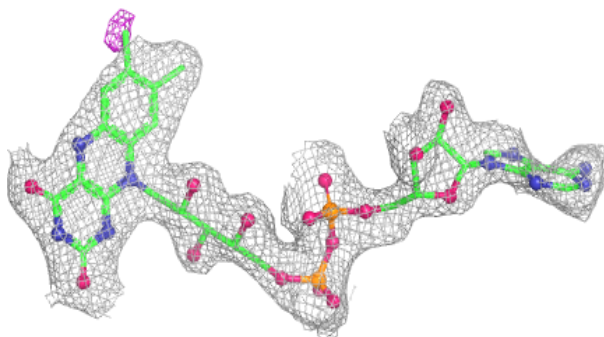
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD A 600:

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

**Electron density around FAD B 600:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.