



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:29 AM UTC

PDB ID : 6WP5 / pdb\_00006wp5  
Title : Pyruvate Kinase M2 mutant-S37D  
Authors : Nandi, S.; Razzaghi, M.; Srivastava, D.; Dey, M.  
Deposited on : 2020-04-26  
Resolution : 2.17 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](mailto: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>.

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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)  
Xtrriage (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

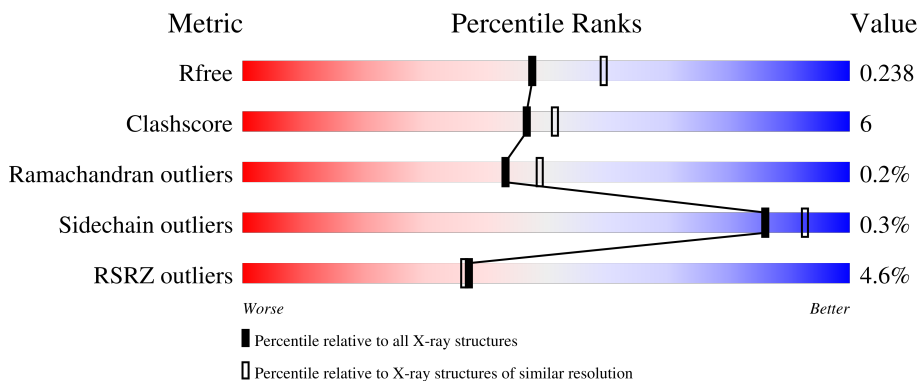
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.17 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	8975 (2.20-2.16)
Clashscore	190562	9786 (2.20-2.16)
Ramachandran outliers	187476	9664 (2.20-2.16)
Sidechain outliers	187428	9664 (2.20-2.16)
RSRZ outliers	180081	8979 (2.20-2.16)

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  $\geq 3$ , 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	550	 2% 80% 12% 7%
1	B	550	 3% 83% 10% 7%
1	C	550	 6% 78% 14% 8%
1	D	550	 5% 65% 9% 26%

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
3	GOL	A	605	-	-	X	-
3	GOL	C	603	-	-	X	-
5	OXL	A	607	-	X	-	-
5	OXL	B	606	-	X	-	-
5	OXL	C	605	-	X	-	-
5	OXL	D	602	-	X	-	-

## 2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 14493 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 Pyruvate kinase PKM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	510	Total 3776	C 2386	N 660	O 706	S 24	0	0	0
1	A	510	Total 3810	C 2408	N 667	O 710	S 25	0	0	0
1	C	507	Total 3641	C 2290	N 645	O 681	S 25	0	0	0
1	D	407	Total 2957	C 1855	N 520	O 560	S 22	0	0	0

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	MET	-	initiating methionine	UNP P14618
B	-17	GLY	-	expression tag	UNP P14618
B	-16	SER	-	expression tag	UNP P14618
B	-15	SER	-	expression tag	UNP P14618
B	-14	HIS	-	expression tag	UNP P14618
B	-13	HIS	-	expression tag	UNP P14618
B	-12	HIS	-	expression tag	UNP P14618
B	-11	HIS	-	expression tag	UNP P14618
B	-10	HIS	-	expression tag	UNP P14618
B	-9	HIS	-	expression tag	UNP P14618
B	-8	SER	-	expression tag	UNP P14618
B	-7	SER	-	expression tag	UNP P14618
B	-6	GLY	-	expression tag	UNP P14618
B	-5	LEU	-	expression tag	UNP P14618
B	-4	VAL	-	expression tag	UNP P14618
B	-3	PRO	-	expression tag	UNP P14618
B	-2	ARG	-	expression tag	UNP P14618
B	-1	GLY	-	expression tag	UNP P14618
B	0	SER	-	expression tag	UNP P14618
B	37	ASP	SER	engineered mutation	UNP P14618
A	-18	MET	-	initiating methionine	UNP P14618

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	GLY	-	expression tag	UNP P14618
A	-16	SER	-	expression tag	UNP P14618
A	-15	SER	-	expression tag	UNP P14618
A	-14	HIS	-	expression tag	UNP P14618
A	-13	HIS	-	expression tag	UNP P14618
A	-12	HIS	-	expression tag	UNP P14618
A	-11	HIS	-	expression tag	UNP P14618
A	-10	HIS	-	expression tag	UNP P14618
A	-9	HIS	-	expression tag	UNP P14618
A	-8	SER	-	expression tag	UNP P14618
A	-7	SER	-	expression tag	UNP P14618
A	-6	GLY	-	expression tag	UNP P14618
A	-5	LEU	-	expression tag	UNP P14618
A	-4	VAL	-	expression tag	UNP P14618
A	-3	PRO	-	expression tag	UNP P14618
A	-2	ARG	-	expression tag	UNP P14618
A	-1	GLY	-	expression tag	UNP P14618
A	0	SER	-	expression tag	UNP P14618
A	37	ASP	SER	engineered mutation	UNP P14618
C	-18	MET	-	initiating methionine	UNP P14618
C	-17	GLY	-	expression tag	UNP P14618
C	-16	SER	-	expression tag	UNP P14618
C	-15	SER	-	expression tag	UNP P14618
C	-14	HIS	-	expression tag	UNP P14618
C	-13	HIS	-	expression tag	UNP P14618
C	-12	HIS	-	expression tag	UNP P14618
C	-11	HIS	-	expression tag	UNP P14618
C	-10	HIS	-	expression tag	UNP P14618
C	-9	HIS	-	expression tag	UNP P14618
C	-8	SER	-	expression tag	UNP P14618
C	-7	SER	-	expression tag	UNP P14618
C	-6	GLY	-	expression tag	UNP P14618
C	-5	LEU	-	expression tag	UNP P14618
C	-4	VAL	-	expression tag	UNP P14618
C	-3	PRO	-	expression tag	UNP P14618
C	-2	ARG	-	expression tag	UNP P14618
C	-1	GLY	-	expression tag	UNP P14618
C	0	SER	-	expression tag	UNP P14618
C	37	ASP	SER	engineered mutation	UNP P14618
D	-18	MET	-	initiating methionine	UNP P14618
D	-17	GLY	-	expression tag	UNP P14618
D	-16	SER	-	expression tag	UNP P14618

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-15	SER	-	expression tag	UNP P14618
D	-14	HIS	-	expression tag	UNP P14618
D	-13	HIS	-	expression tag	UNP P14618
D	-12	HIS	-	expression tag	UNP P14618
D	-11	HIS	-	expression tag	UNP P14618
D	-10	HIS	-	expression tag	UNP P14618
D	-9	HIS	-	expression tag	UNP P14618
D	-8	SER	-	expression tag	UNP P14618
D	-7	SER	-	expression tag	UNP P14618
D	-6	GLY	-	expression tag	UNP P14618
D	-5	LEU	-	expression tag	UNP P14618
D	-4	VAL	-	expression tag	UNP P14618
D	-3	PRO	-	expression tag	UNP P14618
D	-2	ARG	-	expression tag	UNP P14618
D	-1	GLY	-	expression tag	UNP P14618
D	0	SER	-	expression tag	UNP P14618
D	37	ASP	SER	engineered mutation	UNP P14618

- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total K 2 2	0	0
2	A	1	Total K 1 1	0	0
2	C	1	Total K 1 1	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).

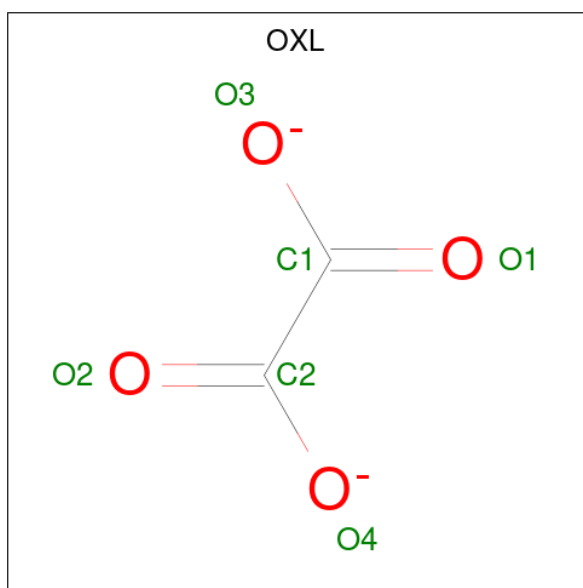


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0
3	D	1	Total C O 6 3 3	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

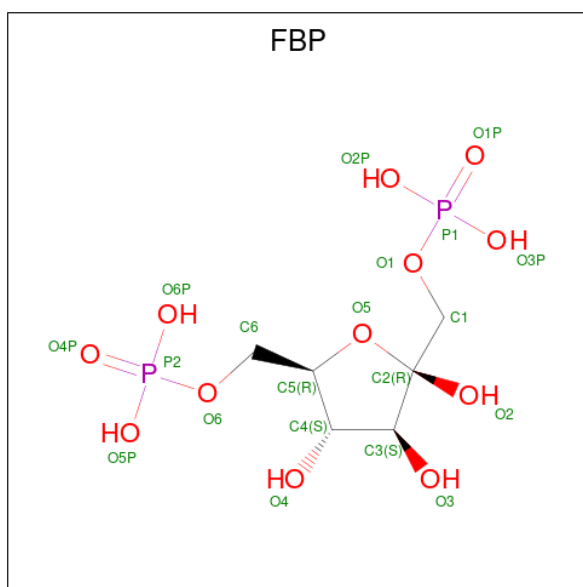
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Mg 1 1	0	0
4	A	1	Total Mg 1 1	0	0
4	C	1	Total Mg 1 1	0	0

- Molecule 5 is OXALATE ION (CCD ID: OXL) (formula:  $C_2O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	2	4		
5	A	1	Total	C	O	0	0
			6	2	4		
5	C	1	Total	C	O	0	0
			6	2	4		
5	D	1	Total	C	O	0	0
			6	2	4		

- Molecule 6 is 1,6-di-O-phosphono-beta-D-fructofuranose (CCD ID: FBP) (formula:  $C_6H_{14}O_{12}P_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total O P 5 4 1	0	0
6	A	1	Total O P 5 4 1	0	0
6	D	1	Total O P 5 4 1	0	0

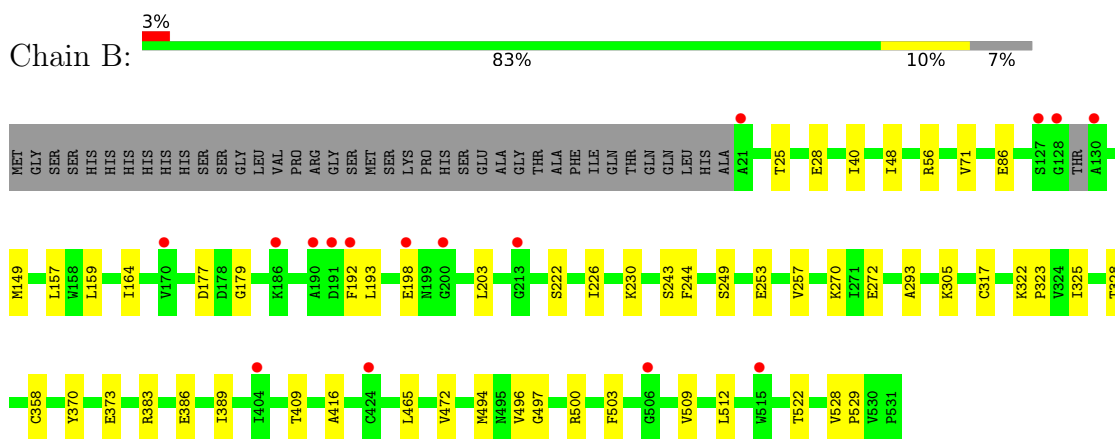
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	67	Total O 67 67	0	0
7	A	75	Total O 75 75	0	0
7	C	29	Total O 29 29	0	0
7	D	38	Total O 38 38	0	0

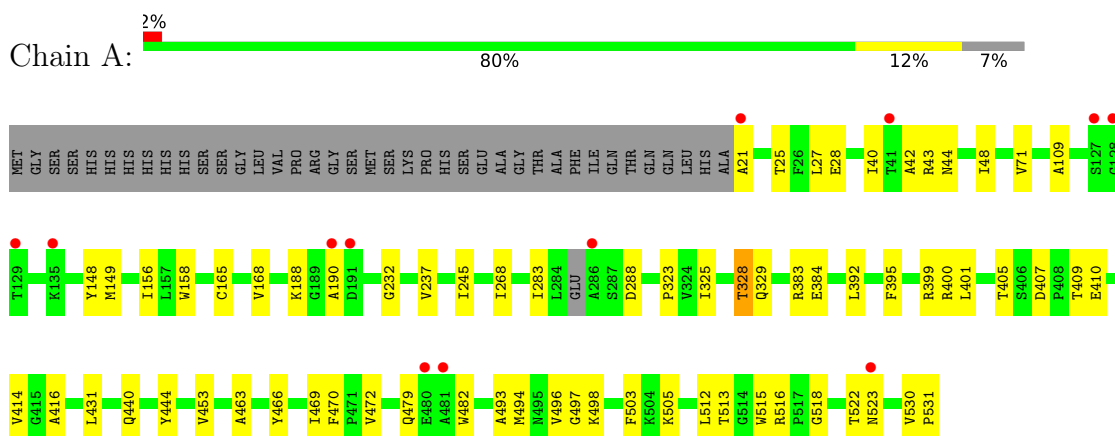
### 3 Residue-property plots i

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.

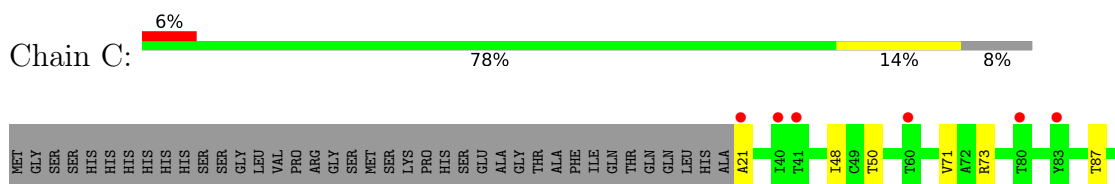
- Molecule 1: Pyruvate kinase PKM



- Molecule 1: Pyruvate kinase PKM



- Molecule 1: Pyruvate kinase PKM





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.62Å 117.57Å 109.97Å 90.00° 112.63° 90.00°	Depositor
Resolution (Å)	19.84 – 2.17 19.84 – 2.17	Depositor EDS
% Data completeness (in resolution range)	99.7 (19.84-2.17) 92.4 (19.84-2.17)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.78 (at 2.17Å)	Xtrriage
Refinement program	PHENIX 1.18.2	Depositor
R, $R_{free}$	0.216 , 0.260 (Not available) , 0.238	Depositor DCC
$R_{free}$ test set	5771 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.8	Xtrriage
Anisotropy	0.732	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 38.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	14493	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.01% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL, MG, OXL, K, FBP

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	0.39	0/3872	0.58	2/5244 (0.0%)
1	B	0.38	0/3837	0.54	1/5203 (0.0%)
1	C	0.47	0/3699	0.58	1/5031 (0.0%)
1	D	0.36	0/3004	0.56	2/4085 (0.0%)
All	All	0.40	0/14412	0.57	6/19563 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	100	SER	CB-CA-C	9.01	128.34	110.42
1	D	22	MET	CB-CA-C	8.48	127.31	110.42
1	B	177	ASP	N-CA-C	7.74	119.97	110.91
1	A	190	ALA	N-CA-C	6.95	125.61	110.80
1	D	362	SER	N-CA-C	6.94	119.93	108.48
1	A	523	ASN	CB-CA-C	5.86	122.09	110.42

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	3810	0	3801	44	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3776	0	3732	29	0
1	C	3641	0	3480	64	0
1	D	2957	0	2835	32	1
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
3	A	24	0	31	7	0
3	B	12	0	16	0	0
3	C	12	0	16	14	0
3	D	6	0	8	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	A	6	0	0	0	0
5	B	6	0	0	0	0
5	C	6	0	0	0	0
5	D	6	0	0	0	0
6	A	5	0	0	0	0
6	B	5	0	0	0	0
6	D	5	0	0	0	0
7	A	75	0	0	1	0
7	B	67	0	0	0	0
7	C	29	0	0	0	0
7	D	38	0	0	1	0
All	All	14493	0	13919	162	1

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 6.

All (162) 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:C:433:LYS:N	3:C:603:GOL:H31	1.39	1.37
1:C:433:LYS:H	3:C:603:GOL:C3	1.47	1.25
1:C:433:LYS:HB3	3:C:603:GOL:H32	1.42	1.01
1:B:244:PHE:N	1:B:272:GLU:OE2	1.98	0.95
1:D:494:MET:HE2	1:D:531:PRO:HD2	1.49	0.94
1:C:432:THR:OG1	3:C:603:GOL:H11	1.69	0.92
1:D:494:MET:HE1	1:D:530:VAL:HG22	1.63	0.80
1:C:433:LYS:HB3	3:C:603:GOL:C3	2.11	0.80
1:C:203:LEU:HD12	1:C:204:GLY:H	1.50	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:494:MET:HE1	1:A:530:VAL:HG22	1.67	0.75
1:A:165:CYS:O	1:A:188:LYS:NZ	2.23	0.72
1:B:472:VAL:HG11	1:B:496:VAL:HG11	1.75	0.69
1:C:142:ILE:HD11	1:C:193:LEU:HD12	1.77	0.67
1:D:395:PHE:CZ	1:D:399:ARG:HD2	2.30	0.66
1:D:466:TYR:HB2	1:D:469:ILE:HD12	1.76	0.66
1:A:328:THR:HG22	1:A:329:GLN:HG3	1.78	0.63
1:C:433:LYS:CB	3:C:603:GOL:H32	2.23	0.62
1:D:50:THR:OG1	1:D:73:ARG:NH1	2.28	0.62
1:A:405:THR:HG22	1:A:407:ASP:H	1.66	0.60
1:B:409:THR:HG23	1:B:522:THR:HB	1.84	0.60
1:B:272:GLU:HG2	1:B:293:ALA:HB3	1.83	0.59
1:A:482:TRP:HH2	3:A:605:GOL:H12	1.69	0.58
1:C:146:ASN:HB3	1:C:149:MET:HE3	1.86	0.57
1:A:494:MET:HE2	1:A:531:PRO:HD2	1.85	0.57
1:D:472:VAL:HG11	1:D:496:VAL:HG11	1.86	0.56
1:C:437:SER:HB3	3:C:603:GOL:O1	2.06	0.56
1:A:42:ALA:HB1	1:A:470:PHE:CZ	2.40	0.56
1:A:472:VAL:HG11	1:A:496:VAL:HG11	1.86	0.56
1:C:21:ALA:HB1	1:C:392:LEU:H	1.70	0.56
1:C:144:LEU:HD13	1:C:162:LYS:HA	1.87	0.56
1:B:370:TYR:HB3	1:B:373:GLU:HB2	1.87	0.55
1:A:25:THR:HB	1:C:397:GLU:CD	2.31	0.55
1:D:21:ALA:N	1:D:392:LEU:H	2.05	0.55
1:C:432:THR:OG1	3:C:603:GOL:C1	2.48	0.55
1:A:466:TYR:HB2	1:A:469:ILE:HD12	1.90	0.54
1:C:433:LYS:HG3	1:C:434:SER:N	2.23	0.54
1:D:81:HIS:HB3	1:D:231:PHE:CE2	2.43	0.54
1:C:425:SER:HB2	1:C:509:VAL:HA	1.90	0.53
1:A:414:VAL:HG12	1:A:444:TYR:CE2	2.43	0.53
1:A:518:GLY:O	3:A:605:GOL:O2	2.25	0.53
1:D:505:LYS:HA	1:D:530:VAL:HG12	1.90	0.53
1:A:505:LYS:HA	1:A:530:VAL:HG12	1.91	0.53
1:B:48:ILE:HG12	1:B:71:VAL:HB	1.90	0.53
1:A:497:GLY:HA3	1:A:503:PHE:CZ	2.44	0.53
1:D:94:ALA:O	1:D:97:SER:HB3	2.09	0.53
1:B:28:GLU:HG3	1:D:401:LEU:HD11	1.90	0.52
1:B:497:GLY:HA3	1:B:503:PHE:CZ	2.44	0.52
1:C:395:PHE:CZ	1:C:399:ARG:HD3	2.44	0.52
1:A:409:THR:HG23	1:A:522:THR:HB	1.90	0.52
1:D:272:GLU:HG2	1:D:293:ALA:HB3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:416:ALA:HB2	1:B:512:LEU:HD21	1.90	0.52
1:D:497:GLY:HA3	1:D:503:PHE:CZ	2.45	0.52
1:D:527:VAL:O	7:D:701:HOH:O	2.19	0.51
1:D:370:TYR:HB3	1:D:373:GLU:HB2	1.91	0.51
1:B:40:ILE:O	1:B:383:ARG:HD2	2.11	0.51
1:D:67:SER:HB3	1:D:376:ARG:HG3	1.92	0.51
1:B:386:GLU:HA	1:B:389:ILE:HD12	1.92	0.51
1:B:528:VAL:HG22	1:B:529:PRO:O	2.12	0.50
1:A:42:ALA:HA	3:A:604:GOL:H32	1.94	0.50
1:A:416:ALA:HB2	1:A:512:LEU:HD21	1.94	0.50
1:C:268:ILE:HD12	1:C:289:GLY:HA3	1.93	0.50
1:D:297:LEU:O	1:D:301:ILE:HG12	2.12	0.50
1:B:192:PHE:O	1:B:193:LEU:HD23	2.12	0.49
1:A:410:GLU:HG3	1:A:440:GLN:HE21	1.78	0.49
1:C:124:ILE:HD11	1:C:203:LEU:HG	1.95	0.49
1:B:179:GLY:O	1:D:339:ARG:NH1	2.46	0.49
1:A:44:ASN:O	3:A:603:GOL:O1	2.25	0.48
1:A:71:VAL:HG22	1:A:109:ALA:HB3	1.95	0.48
1:C:144:LEU:HD11	1:C:164:ILE:HG22	1.94	0.48
1:C:327:ALA:HB1	1:C:360:MET:HE2	1.95	0.48
1:C:395:PHE:O	1:C:399:ARG:HG3	2.13	0.48
1:C:472:VAL:HG11	1:C:496:VAL:HG11	1.96	0.48
1:A:431:LEU:HG	1:A:513:THR:HG22	1.93	0.48
1:C:503:PHE:CD1	1:C:530:VAL:HG21	2.48	0.48
1:C:433:LYS:CA	3:C:603:GOL:H31	2.34	0.48
1:C:433:LYS:CB	3:C:603:GOL:C3	2.87	0.48
1:B:305:LYS:NZ	1:D:384:GLU:OE2	2.31	0.48
1:A:21:ALA:HB1	1:A:392:LEU:H	1.79	0.47
1:A:288:ASP:O	1:A:323:PRO:HD2	2.14	0.47
1:C:48:ILE:HG12	1:C:71:VAL:HB	1.95	0.47
1:A:414:VAL:HG12	1:A:444:TYR:CZ	2.49	0.47
1:B:317:CYS:HB3	1:B:322:LYS:O	2.15	0.47
1:C:258:LEU:HD12	1:C:267:ILE:HD11	1.96	0.47
1:A:453:VAL:HG21	1:A:493:ALA:HB2	1.97	0.47
1:B:325:ILE:HG12	1:B:358:CYS:HB2	1.96	0.47
1:C:50:THR:OG1	1:C:73:ARG:NH1	2.38	0.47
1:C:211:LEU:HB3	1:C:214:ALA:HB3	1.97	0.46
1:C:433:LYS:H	3:C:603:GOL:H31	0.53	0.46
1:A:40:ILE:O	1:A:383:ARG:HD2	2.15	0.46
1:A:48:ILE:HG12	1:A:71:VAL:HB	1.96	0.46
1:D:71:VAL:HG22	1:D:109:ALA:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:294:ARG:HA	1:D:297:LEU:HB3	1.97	0.46
1:D:490:VAL:O	1:D:494:MET:HG2	2.15	0.46
1:A:384:GLU:OE2	1:C:305:LYS:NZ	2.38	0.46
1:C:175:TYR:HB3	1:C:179:GLY:HA2	1.98	0.46
1:A:395:PHE:CZ	1:A:399:ARG:HD3	2.51	0.46
1:C:331:LEU:HD12	1:C:361:LEU:HD21	1.98	0.46
1:C:142:ILE:HA	1:C:157:LEU:O	2.17	0.45
1:B:157:LEU:HD13	1:B:203:LEU:HD21	1.98	0.45
1:B:323:PRO:HB3	1:B:465:LEU:O	2.16	0.45
1:D:429:ILE:HG12	1:D:493:ALA:HB1	1.98	0.45
1:C:169:GLU:HA	1:C:188:LYS:HD2	1.96	0.45
1:B:496:VAL:O	1:B:500:ARG:HG3	2.16	0.45
1:B:159:LEU:HD11	1:B:164:ILE:HD12	1.99	0.45
1:C:323:PRO:HB3	1:C:465:LEU:O	2.16	0.45
1:A:28:GLU:OE2	1:C:319:ARG:NH1	2.50	0.45
1:D:22:MET:O	1:D:23:ALA:C	2.60	0.45
1:A:498:LYS:NZ	1:A:531:PRO:O	2.46	0.44
1:C:417:VAL:HG22	1:C:446:PRO:HB3	2.00	0.44
1:C:510:ILE:HD12	1:C:527:VAL:HG22	1.98	0.44
1:A:463:ALA:HB1	1:A:469:ILE:HG21	1.98	0.44
1:C:189:GLY:HA3	1:C:192:PHE:CZ	2.53	0.44
1:D:81:HIS:CD2	1:D:81:HIS:N	2.85	0.44
1:B:243:SER:HA	1:B:270:LYS:HD3	2.00	0.43
1:C:434:SER:OG	3:C:603:GOL:O2	2.22	0.43
1:C:492:PHE:O	1:C:496:VAL:HG23	2.17	0.43
1:C:512:LEU:HA	1:C:524:THR:O	2.18	0.43
1:A:515:TRP:CE3	1:A:516:ARG:HB2	2.54	0.43
1:C:328:THR:HG22	1:C:329:GLN:HG3	2.00	0.43
1:A:27:LEU:HD23	1:C:401:LEU:HD12	2.01	0.43
1:A:494:MET:CE	1:A:530:VAL:HG13	2.49	0.43
1:C:203:LEU:HD12	1:C:204:GLY:N	2.27	0.43
1:A:401:LEU:HD23	1:A:401:LEU:HA	1.77	0.42
1:C:188:LYS:HE2	1:C:188:LYS:HB3	1.66	0.42
1:D:288:ASP:O	1:D:323:PRO:HD2	2.19	0.42
1:B:494:MET:HE1	1:B:509:VAL:HG21	2.00	0.42
3:A:605:GOL:H11	7:A:724:HOH:O	2.20	0.42
1:D:323:PRO:HA	1:D:357:ASP:OD2	2.18	0.42
1:B:230:LYS:HE2	1:B:257:VAL:HG13	2.01	0.42
1:D:323:PRO:HB3	1:D:465:LEU:O	2.20	0.42
1:A:43:ARG:O	3:A:603:GOL:H31	2.20	0.42
1:A:149:MET:HA	1:A:158:TRP:CG	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:482:TRP:CH2	3:A:605:GOL:H12	2.51	0.42
1:C:288:ASP:O	1:C:323:PRO:HD2	2.20	0.42
1:C:434:SER:CB	3:C:603:GOL:O2	2.67	0.42
1:C:498:LYS:HB2	1:C:498:LYS:HE3	1.83	0.42
1:B:249:SER:O	1:B:253:GLU:HG3	2.20	0.42
1:B:222:SER:O	1:B:226:ILE:HG13	2.20	0.41
1:B:149:MET:HE3	1:B:149:MET:HB2	1.76	0.41
1:A:148:TYR:CE2	1:A:156:ILE:HD13	2.55	0.41
1:C:87:THR:O	1:C:91:VAL:HG23	2.20	0.41
1:B:25:THR:HB	1:D:397:GLU:CD	2.45	0.41
1:A:168:VAL:O	1:A:188:LYS:NZ	2.39	0.41
1:C:188:LYS:HA	1:C:193:LEU:HD23	2.02	0.41
1:C:409:THR:OG1	1:C:521:PHE:HB2	2.21	0.41
1:C:497:GLY:HA3	1:C:503:PHE:CZ	2.56	0.41
1:D:241:PHE:HB3	1:D:270:LYS:HD2	2.03	0.41
1:A:232:GLY:O	1:A:237:VAL:HG22	2.21	0.41
1:C:436:ARG:O	1:C:436:ARG:HD2	2.20	0.41
1:C:510:ILE:HG23	1:C:525:MET:HG3	2.03	0.41
1:D:292:VAL:HG12	1:D:294:ARG:HG2	2.03	0.41
1:C:99:ALA:O	1:C:100:SER:C	2.64	0.41
1:C:399:ARG:HG3	1:C:399:ARG:HH11	1.85	0.41
1:A:245:ILE:HG22	1:A:283:ILE:HD13	2.03	0.40
1:C:175:TYR:HA	1:C:181:ILE:O	2.22	0.40
1:C:210:ASN:HA	3:C:602:GOL:H12	2.03	0.40
1:C:280:PHE:HE1	1:C:290:ILE:HG21	1.86	0.40
1:C:124:ILE:HA	1:C:152:CYS:HB2	2.02	0.40
1:B:56:ARG:NH2	1:B:86:GLU:HB3	2.37	0.40
1:A:268:ILE:HG21	1:A:325:ILE:HD12	2.04	0.40
1:C:176:VAL:HA	1:C:208:GLY:O	2.22	0.40
1:D:50:THR:OG1	1:D:73:ARG:HD3	2.22	0.40

All (1) 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:A:400:ARG:NH2	1:D:22:MET:O[1_554]	2.11	0.09

## 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	506/550 (92%)	493 (97%)	12 (2%)	1 (0%)	43	49
1	B	506/550 (92%)	498 (98%)	7 (1%)	1 (0%)	43	49
1	C	503/550 (92%)	489 (97%)	14 (3%)	0	100	100
1	D	399/550 (72%)	389 (98%)	9 (2%)	1 (0%)	36	39
All	All	1914/2200 (87%)	1869 (98%)	42 (2%)	3 (0%)	43	49

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	328	THR
1	B	328	THR
1	A	328	THR

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	390/452 (86%)	389 (100%)	1 (0%)	86	92
1	B	379/452 (84%)	378 (100%)	1 (0%)	86	92
1	C	347/452 (77%)	345 (99%)	2 (1%)	78	87
1	D	290/452 (64%)	290 (100%)	0	100	100
All	All	1406/1808 (78%)	1402 (100%)	4 (0%)	86	92

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	198	GLU
1	A	479	GLN
1	C	115	LYS
1	C	433	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	184	GLN
1	B	393	GLN
1	A	252	HIS
1	A	458	GLN
1	A	479	GLN
1	D	84	HIS
1	D	393	GLN

### 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](#)

Of 23 ligands modelled in this entry, 7 are monoatomic - leaving 16 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	OXL	C	605	4	5,5,5	2.55	1 (20%)	6,6,6	1.94	2 (33%)
3	GOL	C	602	-	5,5,5	0.83	0	5,5,5	1.20	0
6	FBP	D	603	-	4,4,20	1.42	1 (25%)	6,6,32	1.49	1 (16%)
3	GOL	C	603	-	5,5,5	0.26	0	5,5,5	0.32	0
6	FBP	B	607	-	4,4,20	1.40	1 (25%)	6,6,32	1.49	1 (16%)
3	GOL	A	604	-	5,5,5	1.60	1 (20%)	5,5,5	1.02	0
6	FBP	A	608	-	4,4,20	1.41	1 (25%)	6,6,32	1.49	1 (16%)
5	OXL	A	607	4,2	5,5,5	2.54	1 (20%)	6,6,6	1.94	2 (33%)
3	GOL	B	603	-	5,5,5	0.85	0	5,5,5	1.31	0
3	GOL	A	602	-	5,5,5	2.21	2 (40%)	5,5,5	1.04	0
3	GOL	A	605	-	5,5,5	0.26	0	5,5,5	0.32	0
3	GOL	D	601	-	5,5,5	0.99	0	5,5,5	0.89	0
3	GOL	B	604	-	5,5,5	0.81	0	5,5,5	1.20	0
3	GOL	A	603	-	5,5,5	1.04	0	5,5,5	1.08	1 (20%)
5	OXL	D	602	-	5,5,5	2.55	1 (20%)	6,6,6	1.95	2 (33%)
5	OXL	B	606	4	5,5,5	2.55	1 (20%)	6,6,6	1.95	2 (33%)

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	OXL	C	605	4	-	4/4/4/4	-
3	GOL	C	602	-	-	0/4/4/4	-
3	GOL	C	603	-	-	3/4/4/4	-
3	GOL	A	604	-	-	2/4/4/4	-
5	OXL	A	607	4,2	-	4/4/4/4	-
3	GOL	B	603	-	-	3/4/4/4	-
3	GOL	A	602	-	-	1/4/4/4	-
3	GOL	A	605	-	-	4/4/4/4	-
3	GOL	D	601	-	-	2/4/4/4	-
3	GOL	B	604	-	-	0/4/4/4	-
3	GOL	A	603	-	-	4/4/4/4	-
5	OXL	D	602	-	-	4/4/4/4	-
5	OXL	B	606	4	-	4/4/4/4	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	602	OXL	C2-C1	-4.78	1.46	1.54
5	B	606	OXL	C2-C1	-4.77	1.46	1.54
5	C	605	OXL	C2-C1	-4.76	1.46	1.54
5	A	607	OXL	C2-C1	-4.75	1.46	1.54
3	A	602	GOL	O2-C2	-3.63	1.32	1.43
3	A	602	GOL	C1-C2	3.18	1.63	1.51
3	A	604	GOL	C1-C2	2.84	1.62	1.51
6	D	603	FBP	P2-O6	2.35	1.61	1.54
6	A	608	FBP	P2-O6	2.33	1.61	1.54
6	B	607	FBP	P1-O1	2.32	1.61	1.54

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	605	OXL	O3-C1-C2	3.20	119.03	112.83
5	A	607	OXL	O4-C2-C1	3.19	119.03	112.83
5	B	606	OXL	O3-C1-C2	3.19	119.02	112.83
5	D	602	OXL	O3-C1-C2	3.19	119.01	112.83
5	B	606	OXL	O4-C2-C1	3.18	118.99	112.83
5	D	602	OXL	O4-C2-C1	3.17	118.98	112.83
5	A	607	OXL	O3-C1-C2	3.16	118.96	112.83
5	C	605	OXL	O4-C2-C1	3.15	118.94	112.83
6	A	608	FBP	O6P-P2-O5P	2.09	114.41	107.91
6	B	607	FBP	O3P-P1-O2P	2.07	114.36	107.91
6	D	603	FBP	O6P-P2-O5P	2.07	114.36	107.91
3	A	603	GOL	C3-C2-C1	-2.03	104.36	111.80

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	603	GOL	C1-C2-C3-O3
3	A	603	GOL	C1-C2-C3-O3
3	A	603	GOL	O2-C2-C3-O3
3	A	604	GOL	O1-C1-C2-C3
3	A	605	GOL	O1-C1-C2-C3
3	A	605	GOL	C1-C2-C3-O3
3	A	604	GOL	O1-C1-C2-O2
3	A	602	GOL	O1-C1-C2-C3
3	A	603	GOL	O1-C1-C2-C3
3	C	603	GOL	O1-C1-C2-C3
3	D	601	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	B	603	GOL	O2-C2-C3-O3
3	A	605	GOL	O1-C1-C2-O2
3	A	605	GOL	O2-C2-C3-O3
3	C	603	GOL	O1-C1-C2-O2
3	A	603	GOL	O1-C1-C2-O2
3	B	603	GOL	O1-C1-C2-O2
5	B	606	OXL	O3-C1-C2-O4
5	A	607	OXL	O3-C1-C2-O4
5	C	605	OXL	O3-C1-C2-O4
5	D	602	OXL	O3-C1-C2-O4
5	B	606	OXL	O1-C1-C2-O2
5	A	607	OXL	O1-C1-C2-O2
5	C	605	OXL	O1-C1-C2-O2
5	D	602	OXL	O1-C1-C2-O2
3	D	601	GOL	O1-C1-C2-O2
3	C	603	GOL	C1-C2-C3-O3
5	C	605	OXL	O1-C1-C2-O4
5	C	605	OXL	O3-C1-C2-O2
5	B	606	OXL	O1-C1-C2-O4
5	B	606	OXL	O3-C1-C2-O2
5	A	607	OXL	O1-C1-C2-O4
5	A	607	OXL	O3-C1-C2-O2
5	D	602	OXL	O1-C1-C2-O4
5	D	602	OXL	O3-C1-C2-O2

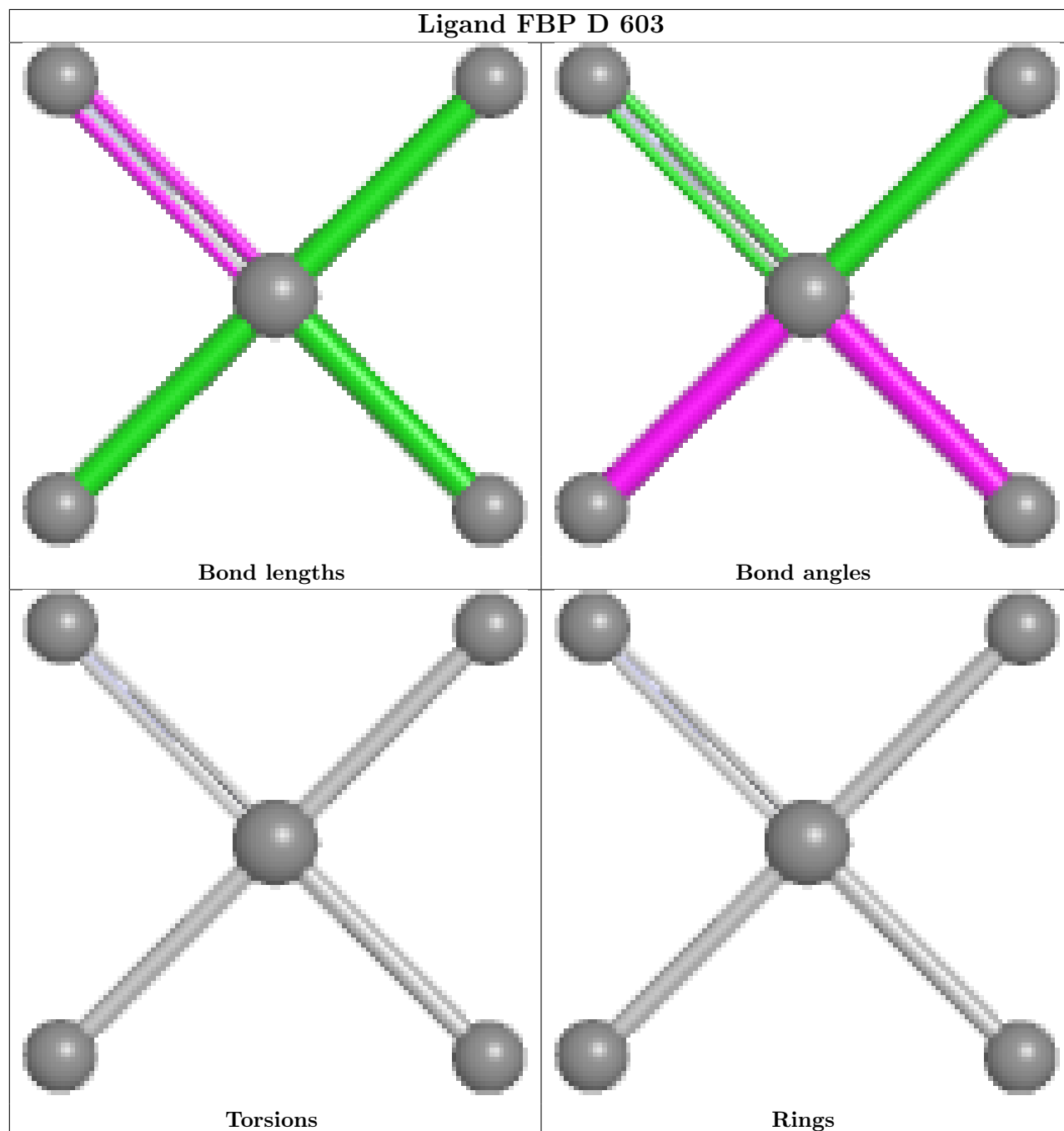
There are no ring outliers.

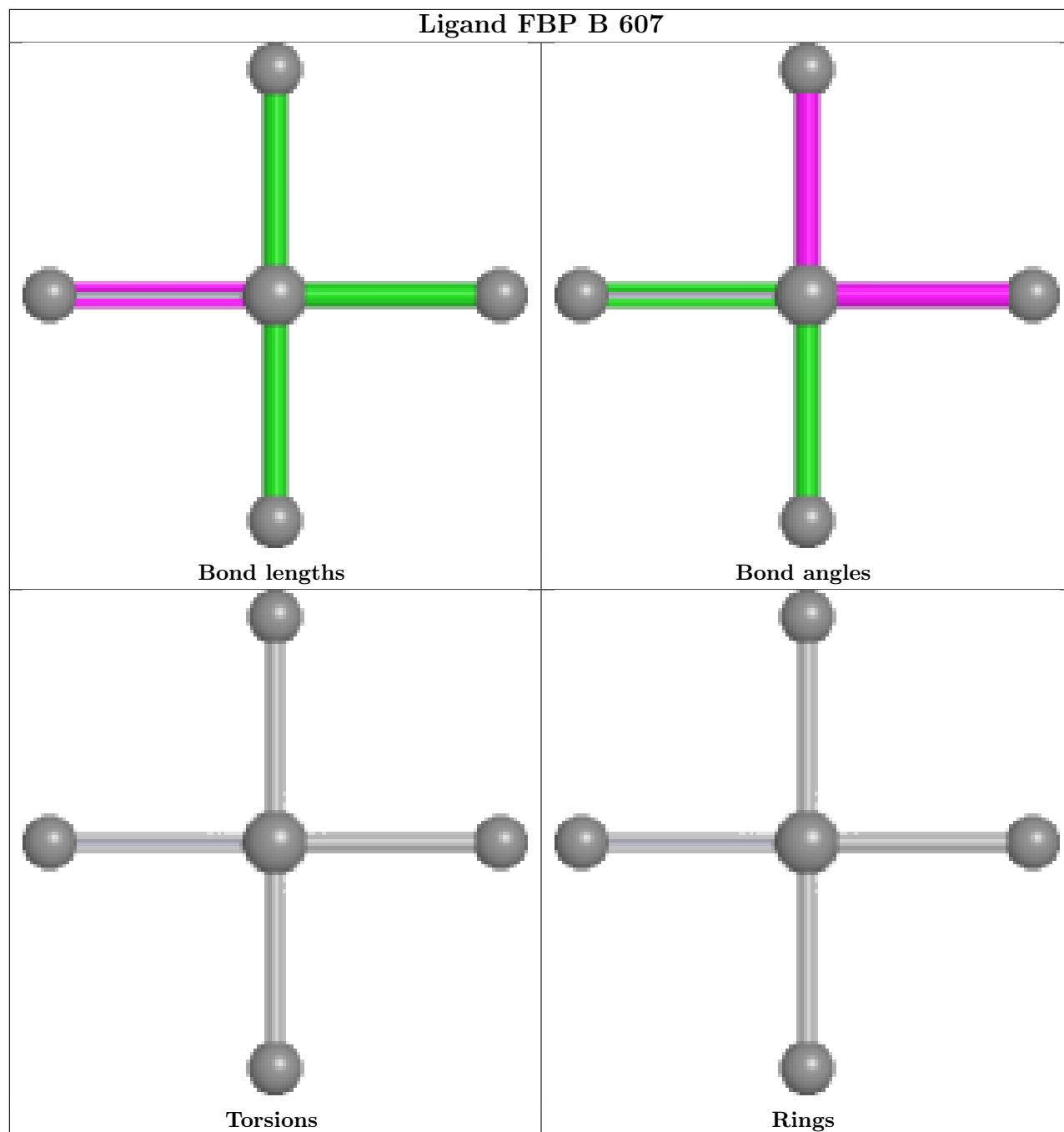
5 monomers are involved in 21 short contacts:

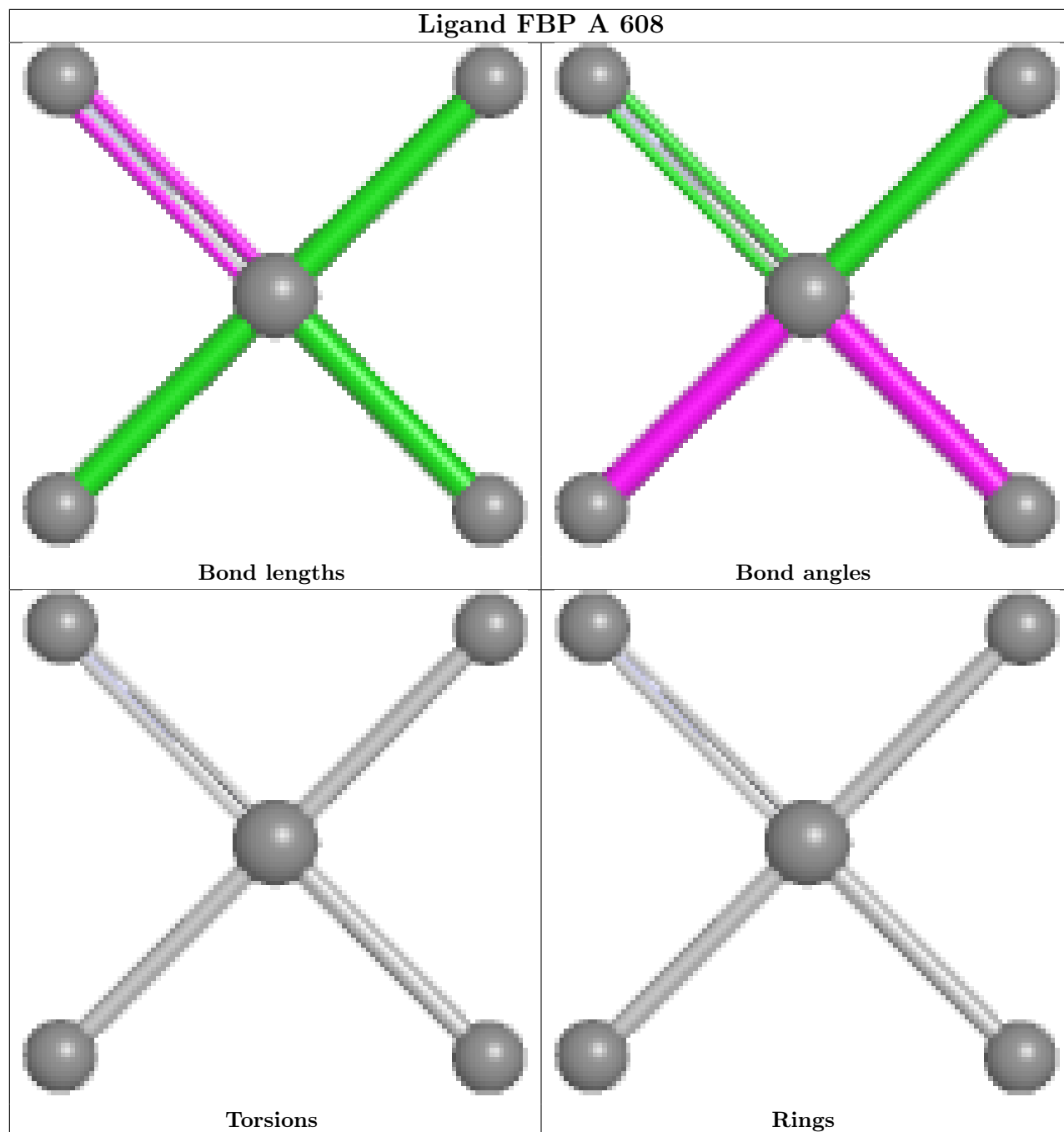
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	602	GOL	1	0
3	C	603	GOL	13	0
3	A	604	GOL	1	0
3	A	605	GOL	4	0
3	A	603	GOL	2	0

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	510/550 (92%)	0.16	12 (2%) 59 59	23, 36, 56, 75	0
1	B	510/550 (92%)	0.20	16 (3%) 51 51	24, 37, 60, 71	0
1	C	507/550 (92%)	0.66	34 (6%) 24 23	25, 44, 67, 79	0
1	D	407/550 (74%)	0.69	27 (6%) 24 23	27, 44, 65, 78	0
All	All	1934/2200 (87%)	0.41	89 (4%) 37 36	23, 40, 64, 79	0

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	21	ALA	5.8
1	B	198	GLU	5.2
1	A	129	THR	4.5
1	C	130	ALA	4.5
1	C	148	TYR	4.2
1	A	21	ALA	4.2
1	A	128	GLY	4.1
1	D	76	PHE	4.0
1	D	119	ILE	3.9
1	D	226	ILE	3.8
1	C	132	VAL	3.8
1	D	520	GLY	3.8
1	A	190	ALA	3.7
1	C	21	ALA	3.7
1	D	245	ILE	3.6
1	C	123	LEU	3.6
1	C	200	GLY	3.6
1	D	21	ALA	3.5
1	C	41	THR	3.5
1	D	262	GLY	3.4
1	C	156	ILE	3.4

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	521	PHE	3.3
1	D	60	THR	3.2
1	C	40	ILE	3.2
1	A	481	ALA	3.2
1	C	124	ILE	3.2
1	B	127	SER	3.2
1	B	191	ASP	3.1
1	A	480	GLU	3.1
1	D	254	VAL	3.1
1	B	190	ALA	3.0
1	C	158	TRP	3.0
1	B	192	PHE	3.0
1	D	251	VAL	2.9
1	C	60	THR	2.8
1	C	191	ASP	2.7
1	D	482	TRP	2.7
1	C	116	GLY	2.7
1	C	133	GLU	2.7
1	B	506	GLY	2.7
1	C	482	TRP	2.6
1	D	220	ALA	2.6
1	D	478	VAL	2.6
1	A	127	SER	2.6
1	D	41	THR	2.6
1	C	117	PRO	2.6
1	A	286	ALA	2.5
1	D	244	PHE	2.5
1	D	273	ASN	2.5
1	A	191	ASP	2.4
1	C	83	TYR	2.4
1	D	219	PRO	2.4
1	C	520	GLY	2.4
1	B	404	ILE	2.4
1	C	203	LEU	2.4
1	A	523	ASN	2.4
1	D	22	MET	2.4
1	C	214	ALA	2.4
1	B	515	TRP	2.4
1	B	130	ALA	2.3
1	B	128	GLY	2.3
1	C	100	SER	2.3
1	D	252	HIS	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	212	PRO	2.3
1	C	515	TRP	2.3
1	C	194	VAL	2.3
1	C	150	GLU	2.2
1	B	200	GLY	2.2
1	D	218	LEU	2.2
1	D	100	SER	2.2
1	A	41	THR	2.2
1	C	91	VAL	2.1
1	C	149	MET	2.1
1	D	78	HIS	2.1
1	C	245	ILE	2.1
1	B	170	VAL	2.1
1	D	229	LEU	2.1
1	D	84	HIS	2.1
1	B	186	LYS	2.1
1	A	135	LYS	2.1
1	D	521	PHE	2.1
1	C	193	LEU	2.0
1	B	424	CYS	2.0
1	B	213	GLY	2.0
1	D	295	GLY	2.0
1	C	131	GLU	2.0
1	C	157	LEU	2.0
1	D	246	ARG	2.0
1	C	80	THR	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 [\(i\)](#)

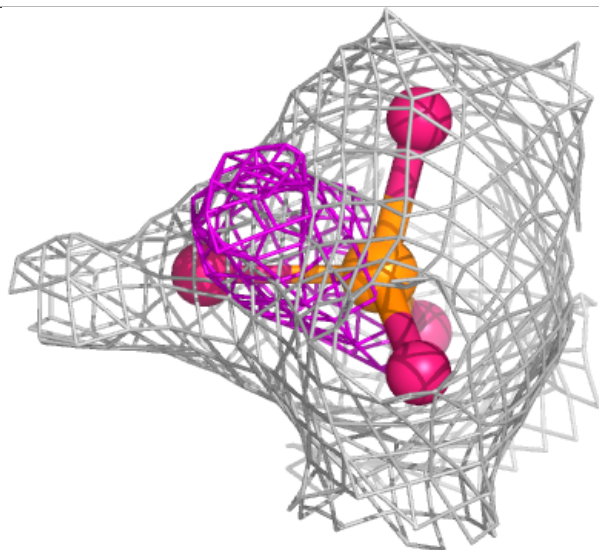
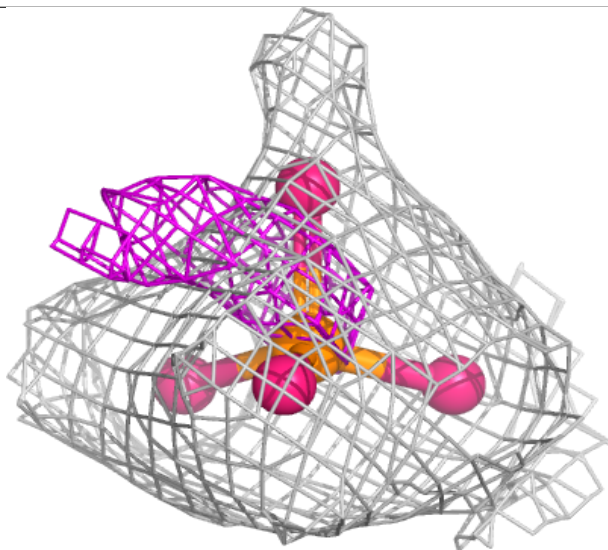
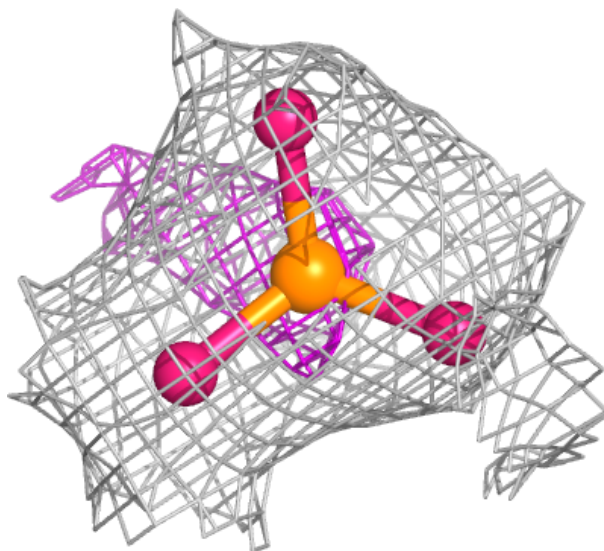
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	K	C	601	1/1	0.64	0.28	77,77,77,77	0
3	GOL	D	601	6/6	0.66	0.17	55,65,67,69	0
5	OXL	D	602	6/6	0.74	0.14	46,48,52,54	0
2	K	B	602	1/1	0.78	0.21	59,59,59,59	0
3	GOL	A	605	6/6	0.78	0.19	41,48,51,57	0
3	GOL	C	603	6/6	0.80	0.12	40,48,50,53	0
3	GOL	C	602	6/6	0.83	0.14	52,56,57,59	0
3	GOL	B	603	6/6	0.83	0.13	36,38,43,44	0
3	GOL	A	604	6/6	0.83	0.17	37,39,46,50	0
2	K	A	601	1/1	0.83	0.29	67,67,67,67	0
3	GOL	A	603	6/6	0.86	0.14	34,36,40,44	0
6	FBP	D	603	5/20	0.86	0.14	45,54,57,58	0
5	OXL	C	605	6/6	0.89	0.10	41,46,46,47	0
3	GOL	A	602	6/6	0.89	0.14	32,37,42,46	0
4	MG	C	604	1/1	0.89	0.12	49,49,49,49	0
3	GOL	B	604	6/6	0.92	0.10	43,45,48,49	0
5	OXL	B	606	6/6	0.95	0.07	29,33,36,38	0
2	K	B	601	1/1	0.95	0.06	41,41,41,41	0
5	OXL	A	607	6/6	0.96	0.06	30,32,36,37	0
6	FBP	B	607	5/20	0.96	0.07	39,40,43,44	0
4	MG	B	605	1/1	0.96	0.05	33,33,33,33	0
6	FBP	A	608	5/20	0.97	0.09	37,37,39,41	0
4	MG	A	606	1/1	0.98	0.05	36,36,36,36	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.

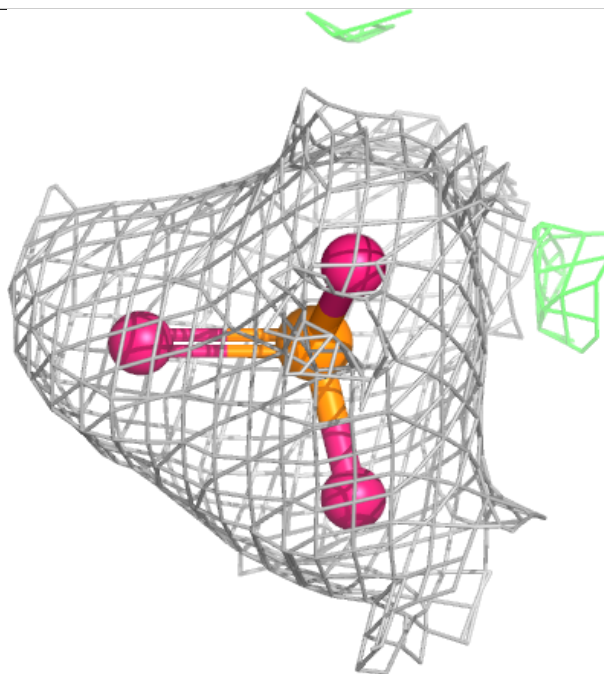
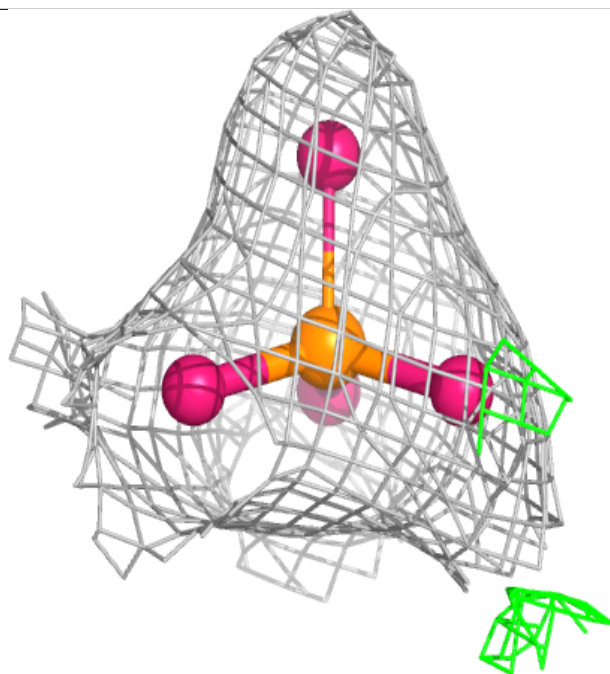
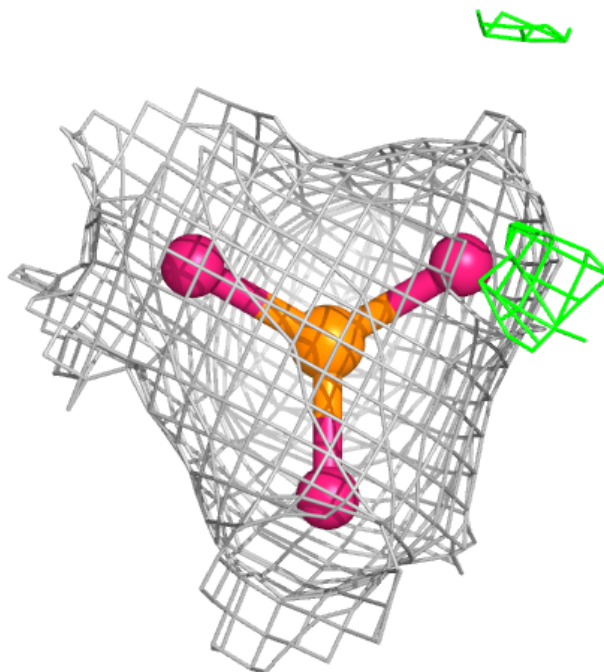
**Electron density around FBP D 603:**

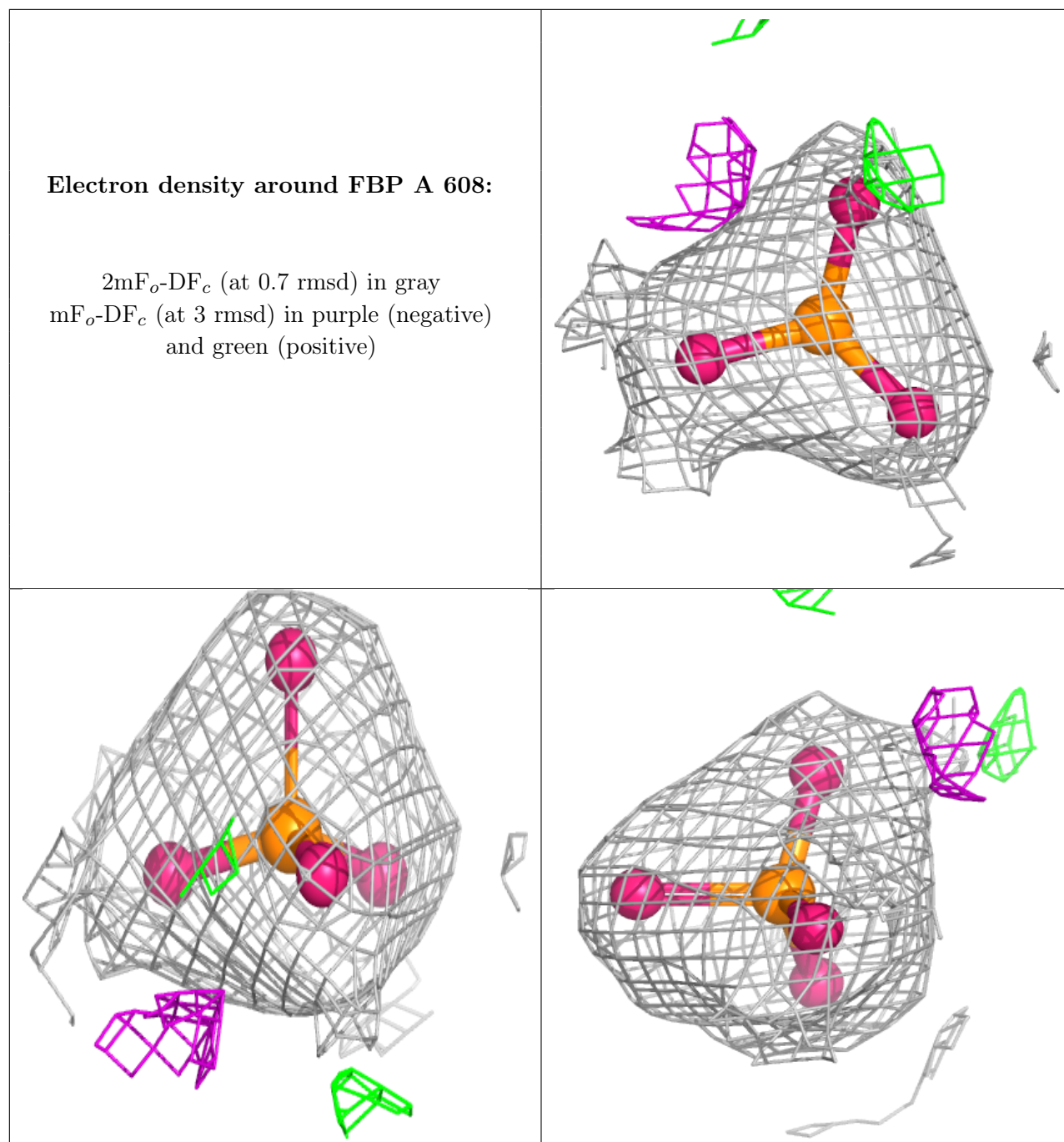
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around FBP B 607:**

$2mF_o-DF_c$  (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.