



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 4, 2026 – 05:02 AM EDT

PDB ID : 4YDD / pdb_00004ydd
Title : Crystal structure of the perchlorate reductase PcrAB from *Azospira suillum*
PS
Authors : Tsai, C.-L.; Youngblut, M.D.; Tainer, J.A.
Deposited on : 2015-02-21
Resolution : 1.86 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

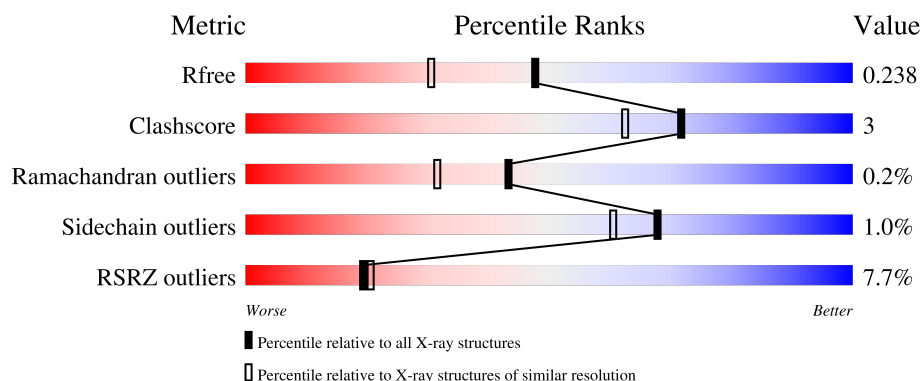
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.86 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	899	93%6%
1	C	899	6% 91%7%
1	E	899	9% 91%8%
2	B	333	92%6%
2	D	333	39% 87%11%

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Mol	Chain	Length	Quality of chain
2	F	333	% 92% 7%

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 61265 atoms, of which 28814 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.

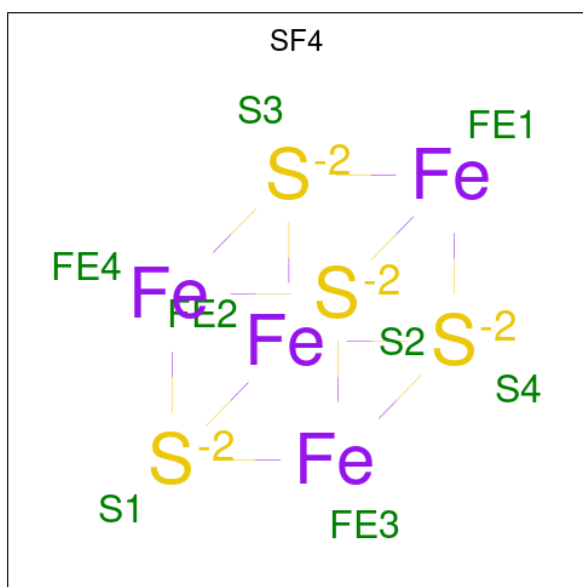
- Molecule 1 is a protein called DMSO reductase family type II enzyme, molybdopterin subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	895	Total	C	H	N	O	S	0	3	0
			14201	4589	7016	1247	1311	38			
1	C	892	Total	C	H	N	O	S	0	0	0
			14110	4563	6962	1240	1307	38			
1	E	892	Total	C	H	N	O	S	0	2	0
			14140	4575	6977	1240	1310	38			

- Molecule 2 is a protein called DMSO reductase family type II enzyme, iron-sulfur subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	329	Total	C	H	N	O	S	0	0	0
			5098	1627	2534	447	465	25			
2	D	328	Total	C	H	N	O	S	0	0	0
			5081	1622	2525	446	464	24			
2	F	328	Total	C	H	N	O	S	0	1	0
			5095	1626	2533	447	465	24			

- Molecule 3 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).

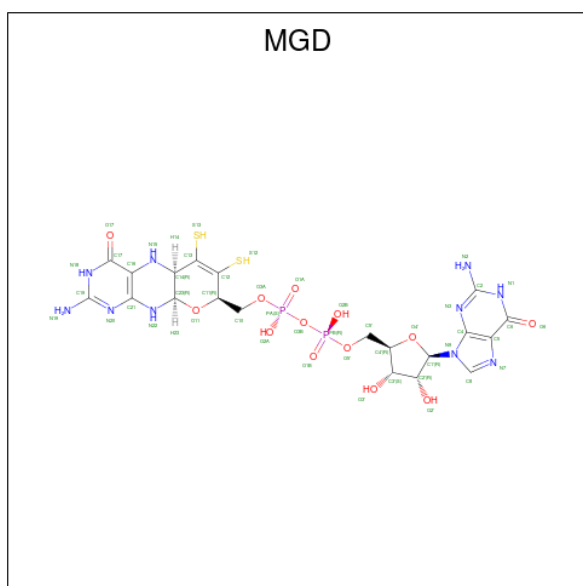


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	S	0	0
			8	4	4		
3	B	1	Total	Fe	S	0	0
			8	4	4		
3	B	1	Total	Fe	S	0	0
			8	4	4		
3	B	1	Total	Fe	S	0	0
			8	4	4		
3	C	1	Total	Fe	S	0	0
			8	4	4		
3	D	1	Total	Fe	S	0	0
			8	4	4		
3	D	1	Total	Fe	S	0	0
			8	4	4		
3	D	1	Total	Fe	S	0	0
			8	4	4		
3	E	1	Total	Fe	S	0	0
			8	4	4		
3	F	1	Total	Fe	S	0	0
			8	4	4		
3	F	1	Total	Fe	S	0	0
			8	4	4		
3	F	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 4 is MOLYBDENUM ATOM (CCD ID: MO) (formula: Mo).

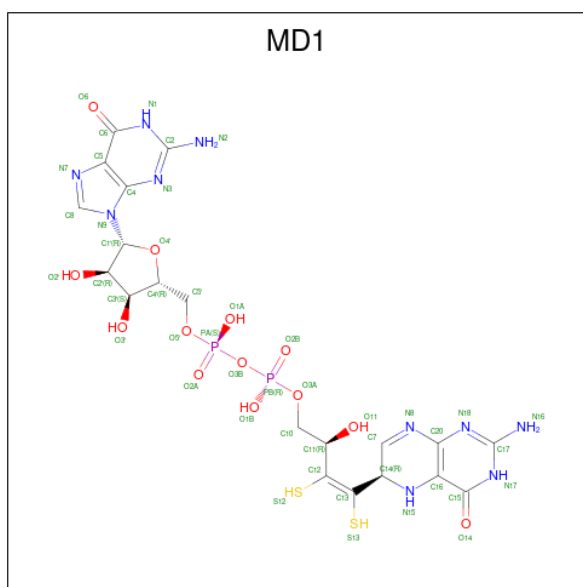
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mo 1 1	0	0
4	C	1	Total Mo 1 1	0	0
4	E	1	Total Mo 1 1	0	0

- Molecule 5 is 2-AMINO-5,6-DIMERCAPTO-7-METHYL-3,7,8A,9-TETRAHYDRO-8-OXA-1,3,9,10-TETRAAZA-ANTHRACEN-4-ONE GUANOSINE DINUCLEOTIDE (CCD ID: MGD) (formula: $C_{20}H_{26}N_{10}O_{13}P_2S_2$).



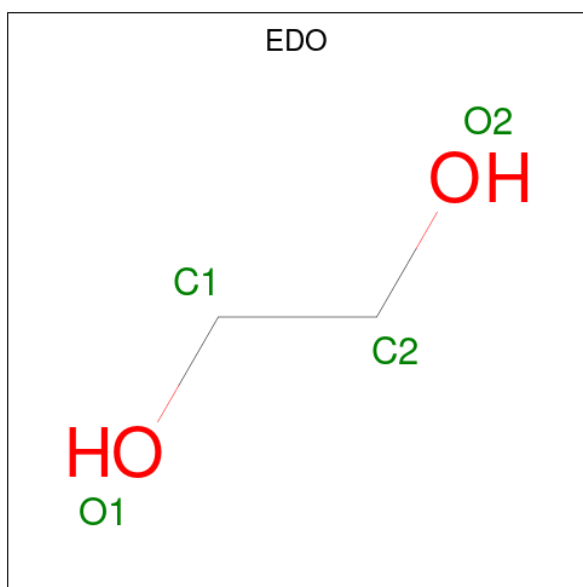
Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	P	S	0	0
			70	20	23	10	13	2	2		
5	C	1	Total	C	H	N	O	P	S	0	0
			70	20	23	10	13	2	2		
5	E	1	Total	C	H	N	O	P	S	0	0
			70	20	23	10	13	2	2		

- Molecule 6 is PHOSPHORIC ACID 4-(2-AMINO-4-OXO-3,4,5,6,-TETRAHYDRO-PTE RIDIN-6-YL)-2-HYDROXY-3,4-DIMERCAPTO-BUT-3-EN-YL ESTER GUANYLATE ESTER (CCD ID: MD1) (formula: $C_{20}H_{26}N_{10}O_{13}P_2S_2$).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
6	A	1	Total 71	C 20	H 24	N 10	O 13	P 2	S 2	0	0
6	C	1	Total 71	C 20	H 24	N 10	O 13	P 2	S 2	0	0
6	E	1	Total 71	C 20	H 24	N 10	O 13	P 2	S 2	0	0

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



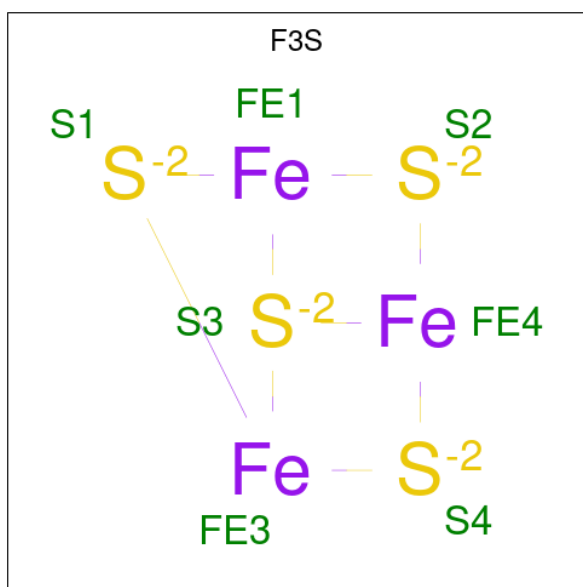
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	H	O	0	0
			10	2	6	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	H	O	0	0
			10	2	6	2		
7	A	1	Total	C	H	O	0	0
			10	2	6	2		
7	A	1	Total	C	H	O	0	0
			10	2	6	2		
7	A	1	Total	C	H	O	0	0
			10	2	6	2		
7	A	1	Total	C	H	O	0	0
			10	2	6	2		
7	B	1	Total	C	H	O	0	0
			10	2	6	2		
7	B	1	Total	C	H	O	0	0
			10	2	6	2		
7	C	1	Total	C	H	O	0	0
			10	2	6	2		
7	C	1	Total	C	H	O	0	0
			10	2	6	2		
7	D	1	Total	C	H	O	0	0
			10	2	6	2		
7	E	1	Total	C	H	O	0	0
			10	2	6	2		
7	E	1	Total	C	H	O	0	0
			10	2	6	2		
7	E	1	Total	C	H	O	0	0
			10	2	6	2		
7	E	1	Total	C	H	O	0	0
			10	2	6	2		
7	E	1	Total	C	H	O	0	0
			10	2	6	2		
7	F	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 8 is FE3-S4 CLUSTER (CCD ID: F3S) (formula: Fe₃S₄).

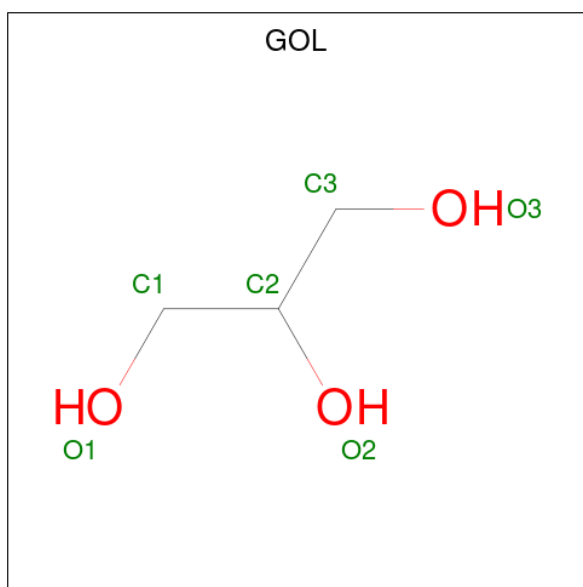


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	Fe	S	0	0
			7	3	4		
8	D	1	Total	Fe	S	0	0
			7	3	4		
8	F	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 9 is SODIUM ION (CCD ID: NA) (formula: Na).

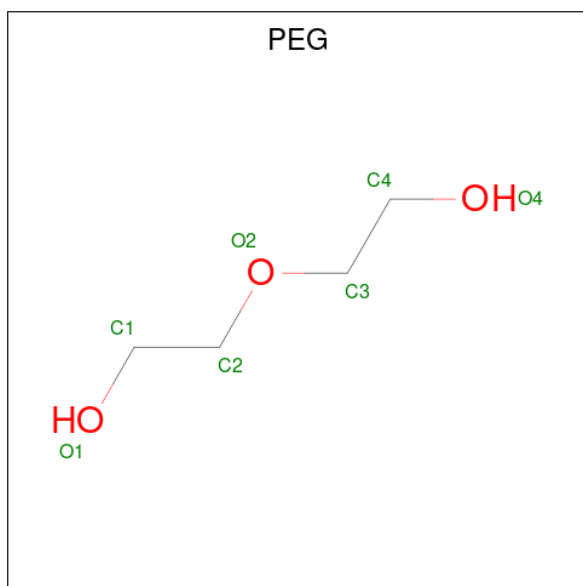
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	1	Total	Na	0	0
			1	1		

- Molecule 10 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	C	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 11 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	C	1	Total	C	H	O	0	0
			17	4	10	3		

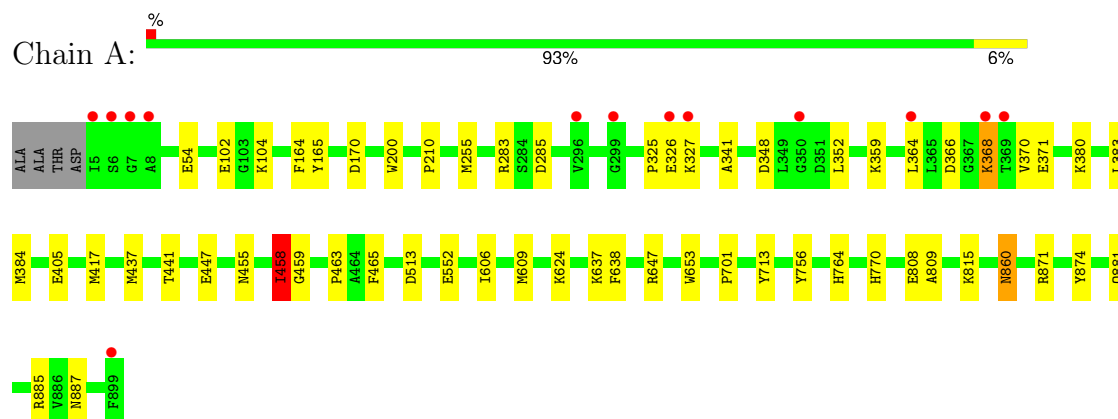
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	738	Total 738	O 738	0	0
12	B	357	Total 358	O 358	0	1
12	C	575	Total 576	O 576	0	1
12	D	140	Total 141	O 141	0	1
12	E	681	Total 681	O 681	0	0
12	F	291	Total 291	O 291	0	0

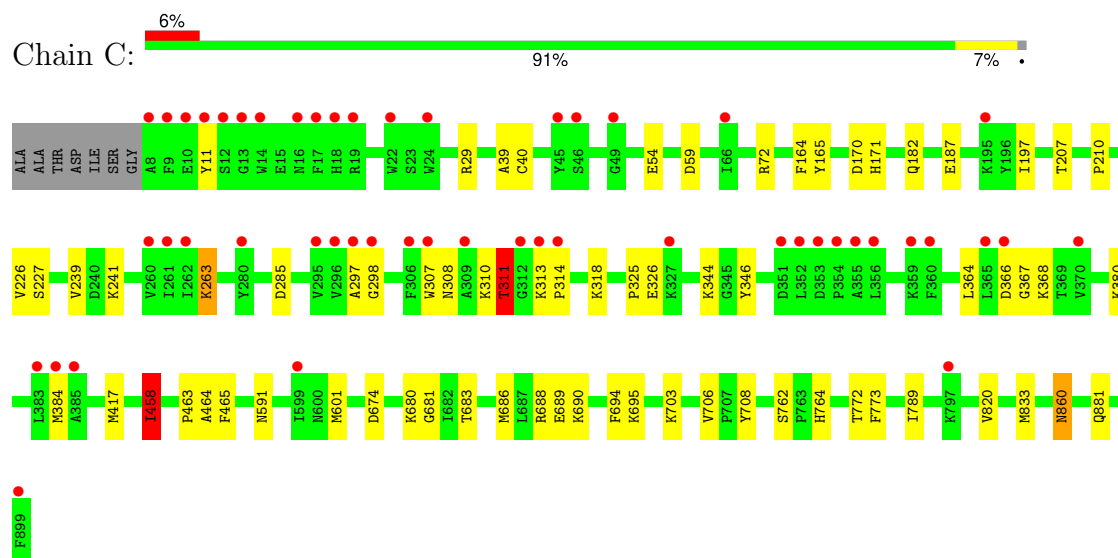
3 Residue-property plots [i](#)

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.

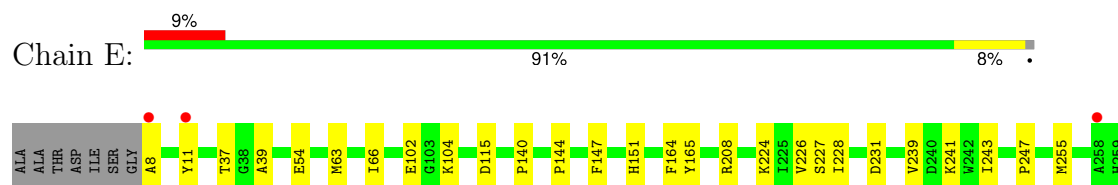
- Molecule 1: DMSO reductase family type II enzyme, molybdopterin subunit

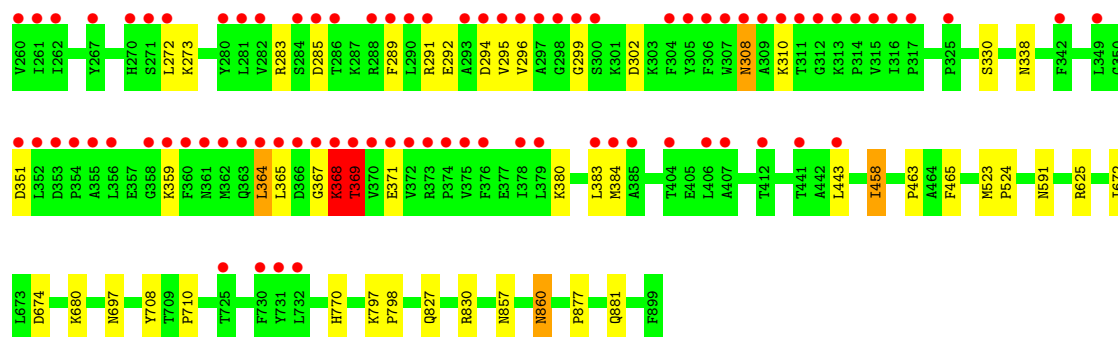


- Molecule 1: DMSO reductase family type II enzyme, molybdopterin subunit



- Molecule 1: DMSO reductase family type II enzyme, molybdopterin subunit

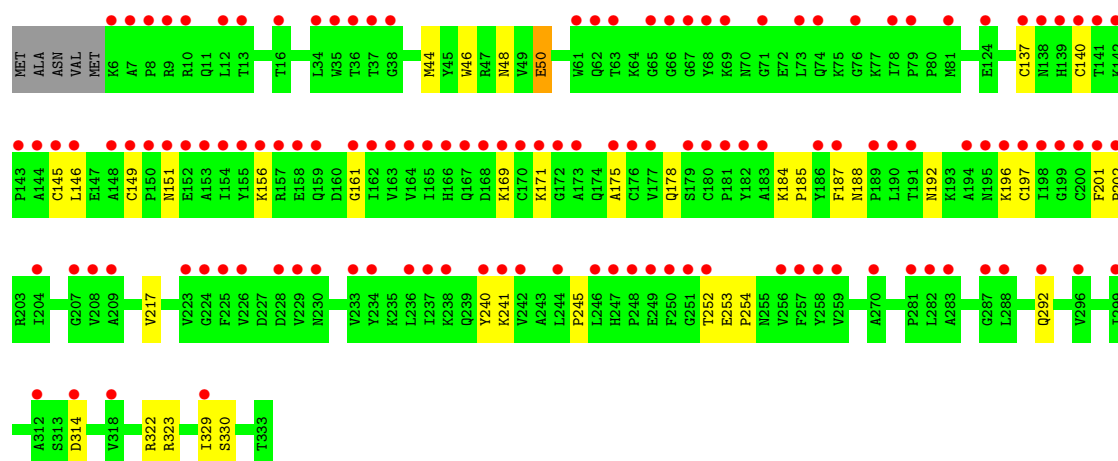
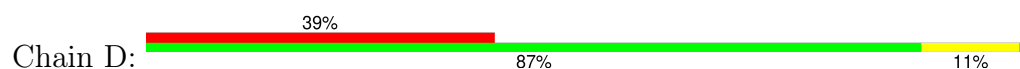




- Molecule 2: DMSO reductase family type II enzyme, iron-sulfur subunit



- Molecule 2: DMSO reductase family type II enzyme, iron-sulfur subunit



- Molecule 2: DMSO reductase family type II enzyme, iron-sulfur subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	132.83Å 175.50Å 193.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.35 – 1.86 48.35 – 1.86	Depositor EDS
% Data completeness (in resolution range)	98.9 (48.35-1.86) 98.9 (48.35-1.86)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 1.86Å)	Xtriage
Refinement program	PHENIX dev_2299	Depositor
R, R_{free}	0.201 , 0.237 0.203 , 0.238	Depositor DCC
R_{free} test set	18608 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	23.1	Xtriage
Anisotropy	0.444	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 35.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	61265	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.04% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PEG, SF4, F3S, MGD, MD1, NA, MO, EDO

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	0.46	0/7403	0.61	0/10045
1	C	0.41	0/7357	0.59	0/9984
1	E	0.45	0/7379	0.63	0/10014
2	B	0.51	0/2632	0.65	0/3567
2	D	0.34	0/2624	0.59	0/3557
2	F	0.45	0/2633	0.62	0/3569
All	All	0.44	0/30028	0.61	0/40736

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7185	7016	7016	34	0
1	C	7148	6962	6960	45	0
1	E	7163	6977	6975	53	0
2	B	2564	2534	2534	14	0
2	D	2556	2525	2525	27	0
2	F	2562	2533	2533	14	0
3	A	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	24	0	0	0	0
3	C	8	0	0	0	0
3	D	24	0	0	1	0
3	E	8	0	0	0	0
3	F	24	0	0	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
5	A	47	23	21	0	0
5	C	47	23	21	1	0
5	E	47	23	21	1	0
6	A	47	24	22	5	0
6	C	47	24	21	4	0
6	E	47	24	21	4	0
7	A	24	36	36	1	0
7	B	8	12	12	0	0
7	C	8	12	12	0	0
7	D	4	6	6	0	0
7	E	24	36	36	0	0
7	F	4	6	6	0	0
8	B	7	0	0	0	0
8	D	7	0	0	1	0
8	F	7	0	0	0	0
9	B	1	0	0	0	0
10	C	6	8	8	0	0
11	C	7	10	10	1	0
12	A	738	0	0	7	1
12	B	358	0	0	2	2
12	C	576	0	0	3	0
12	D	141	0	0	0	0
12	E	681	0	0	7	2
12	F	291	0	0	2	1
All	All	32451	28814	28796	192	3

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 3.

The worst 5 of 192 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:369:THR:O	12:E:1101:HOH:O	2.01	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:231:ASP:OD1	1:E:830:ARG:NH2	2.21	0.74
1:A:405:GLU:OE2	12:A:1101:HOH:O	2.06	0.73
1:E:368:LYS:O	1:E:369:THR:OG1	2.05	0.73
1:C:680:LYS:O	1:C:695:LYS:HE3	1.90	0.72

All (3) 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)
12:B:791:HOH:O	12:E:1471:HOH:O[4_477]	2.04	0.16
12:B:808:HOH:O	12:E:1633:HOH:O[4_477]	2.06	0.14
12:A:1703:HOH:O	12:F:578:HOH:O[2_874]	2.08	0.12

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	896/899 (100%)	864 (96%)	30 (3%)	2 (0%)	43	31
1	C	890/899 (99%)	851 (96%)	36 (4%)	3 (0%)	36	25
1	E	892/899 (99%)	854 (96%)	34 (4%)	4 (0%)	30	17
2	B	327/333 (98%)	316 (97%)	11 (3%)	0	100	100
2	D	326/333 (98%)	314 (96%)	12 (4%)	0	100	100
2	F	327/333 (98%)	316 (97%)	11 (3%)	0	100	100
All	All	3658/3696 (99%)	3515 (96%)	134 (4%)	9 (0%)	43	31

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	311	THR
1	C	458	ILE

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Mol	Chain	Res	Type
1	E	369	THR
1	A	458	ILE
1	E	368	LYS

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	769/768 (100%)	763 (99%)	6 (1%)	73	67
1	C	764/768 (100%)	755 (99%)	9 (1%)	63	55
1	E	766/768 (100%)	757 (99%)	9 (1%)	63	55
2	B	278/281 (99%)	275 (99%)	3 (1%)	65	57
2	D	277/281 (99%)	275 (99%)	2 (1%)	76	70
2	F	278/281 (99%)	275 (99%)	3 (1%)	65	57
All	All	3132/3147 (100%)	3100 (99%)	32 (1%)	68	60

5 of 32 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	881	GLN
2	F	50	GLU
1	C	458	ILE
1	C	311	THR
2	F	72	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	522	ASN
1	E	617	GLN
2	F	151	ASN
1	E	654	ASN
1	C	553	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 45 ligands modelled in this entry, 4 are monoatomic - leaving 41 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$
7	EDO	C	1005	-	3,3,3	0.85	0	2,2,2	0.58	0
6	MD1	E	1004	4	47,51,51	3.08	17 (36%)	56,78,78	1.67	12 (21%)
10	GOL	C	1006	-	5,5,5	0.49	0	5,5,5	0.42	0
7	EDO	A	1010	-	3,3,3	0.63	0	2,2,2	0.62	0
5	MGD	A	1003	4	47,52,52	4.44	28 (59%)	58,81,81	2.33	14 (24%)
5	MGD	E	1003	4	47,52,52	4.63	27 (57%)	58,81,81	2.27	19 (32%)
3	SF4	B	402	2	0,12,12	-	-	-	-	-
7	EDO	A	1007	-	3,3,3	0.46	0	2,2,2	0.40	0
7	EDO	E	1010	-	3,3,3	0.64	0	2,2,2	0.11	0
7	EDO	E	1005	-	3,3,3	0.72	0	2,2,2	0.50	0
7	EDO	E	1008	-	3,3,3	0.29	0	2,2,2	0.61	0
7	EDO	A	1006	-	3,3,3	0.41	0	2,2,2	0.50	0
3	SF4	E	1001	1	0,12,12	-	-	-	-	-
7	EDO	D	405	-	3,3,3	0.47	0	2,2,2	0.42	0
3	SF4	F	402	2	0,12,12	-	-	-	-	-
7	EDO	A	1005	-	3,3,3	0.46	0	2,2,2	0.14	0
6	MD1	C	1004	4	47,51,51	3.25	17 (36%)	56,78,78	1.58	11 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	F3S	F	401	2	0,9,9	-	-	-		
7	EDO	B	405	-	3,3,3	0.44	0	2,2,2	0.93	0
7	EDO	A	1008	-	3,3,3	0.39	0	2,2,2	0.69	0
3	SF4	D	404	2	0,12,12	-	-	-		
11	PEG	C	1008	-	6,6,6	0.59	0	5,5,5	0.57	0
3	SF4	C	1001	1	0,12,12	-	-	-		
7	EDO	A	1009	-	3,3,3	0.50	0	2,2,2	0.63	0
3	SF4	F	403	2	0,12,12	-	-	-		
7	EDO	E	1006	-	3,3,3	0.99	0	2,2,2	0.37	0
7	EDO	F	405	-	3,3,3	0.30	0	2,2,2	0.55	0
8	F3S	B	401	2	0,9,9	-	-	-		
5	MGD	C	1003	4	47,52,52	4.64	28 (59%)	58,81,81	2.43	19 (32%)
7	EDO	C	1007	-	3,3,3	0.49	0	2,2,2	0.41	0
3	SF4	B	404	2	0,12,12	-	-	-		
3	SF4	D	403	2	0,12,12	-	-	-		
6	MD1	A	1004	4	47,51,51	3.16	16 (34%)	56,78,78	1.61	13 (23%)
3	SF4	F	404	2	0,12,12	-	-	-		
3	SF4	D	402	2	0,12,12	-	-	-		
7	EDO	B	406	-	3,3,3	0.72	0	2,2,2	0.32	0
7	EDO	E	1007	-	3,3,3	0.63	0	2,2,2	0.25	0
8	F3S	D	401	2	0,9,9	-	-	-		
3	SF4	B	403	2	0,12,12	-	-	-		
7	EDO	E	1009	-	3,3,3	0.55	0	2,2,2	0.46	0
3	SF4	A	1001	1	0,12,12	-	-	-		

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
7	EDO	C	1005	-	-	1/1/1/1	-
6	MD1	E	1004	4	-	1/22/59/59	0/5/5/5
10	GOL	C	1006	-	-	2/4/4/4	-
7	EDO	A	1010	-	-	0/1/1/1	-
5	MGD	A	1003	4	-	2/22/66/66	0/6/6/6
5	MGD	E	1003	4	-	2/22/66/66	0/6/6/6
3	SF4	B	402	2	-	-	0/6/5/5
7	EDO	A	1007	-	-	1/1/1/1	-
7	EDO	E	1010	-	-	1/1/1/1	-
7	EDO	E	1005	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	EDO	E	1008	-	-	1/1/1/1	-
7	EDO	A	1006	-	-	0/1/1/1	-
3	SF4	A	1001	1	-	-	0/6/5/5
3	SF4	E	1001	1	-	-	0/6/5/5
7	EDO	D	405	-	-	0/1/1/1	-
3	SF4	F	402	2	-	-	0/6/5/5
7	EDO	A	1005	-	-	0/1/1/1	-
6	MD1	C	1004	4	-	6/22/59/59	0/5/5/5
8	F3S	F	401	2	-	-	0/3/3/3
7	EDO	B	405	-	-	0/1/1/1	-
7	EDO	A	1008	-	-	1/1/1/1	-
3	SF4	D	404	2	-	-	0/6/5/5
11	PEG	C	1008	-	-	3/4/4/4	-
3	SF4	C	1001	1	-	-	0/6/5/5
7	EDO	A	1009	-	-	0/1/1/1	-
3	SF4	F	403	2	-	-	0/6/5/5
7	EDO	E	1006	-	-	0/1/1/1	-
7	EDO	F	405	-	-	0/1/1/1	-
8	F3S	B	401	2	-	-	0/3/3/3
7	EDO	C	1007	-	-	1/1/1/1	-
5	MGD	C	1003	4	-	2/22/66/66	0/6/6/6
3	SF4	B	404	2	-	-	0/6/5/5
3	SF4	D	403	2	-	-	0/6/5/5
6	MD1	A	1004	4	-	2/22/59/59	0/5/5/5
7	EDO	B	406	-	-	0/1/1/1	-
3	SF4	D	402	2	-	-	0/6/5/5
3	SF4	F	404	2	-	-	0/6/5/5
8	F3S	D	401	2	-	-	0/3/3/3
3	SF4	B	403	2	-	-	0/6/5/5
7	EDO	E	1009	-	-	1/1/1/1	-
7	EDO	E	1007	-	-	1/1/1/1	-

The worst 5 of 133 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1003	MGD	C16-C21	13.16	1.61	1.38
5	E	1003	MGD	C16-C21	12.33	1.59	1.38
6	C	1004	MD1	C7-N8	12.30	1.43	1.28
5	C	1003	MGD	C16-C21	12.12	1.59	1.38
6	A	1004	MD1	C7-N8	12.03	1.43	1.28

The worst 5 of 88 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1003	MGD	O11-C23-N22	-12.22	97.52	108.61
5	C	1003	MGD	O11-C23-N22	-11.76	97.93	108.61
5	E	1003	MGD	O11-C23-N22	-10.08	99.46	108.61
6	C	1004	MD1	C5-C4-N3	-5.15	120.20	128.39
5	C	1003	MGD	C23-C14-N15	5.13	112.91	107.87

There are no chirality outliers.

5 of 28 torsion outliers are listed below:

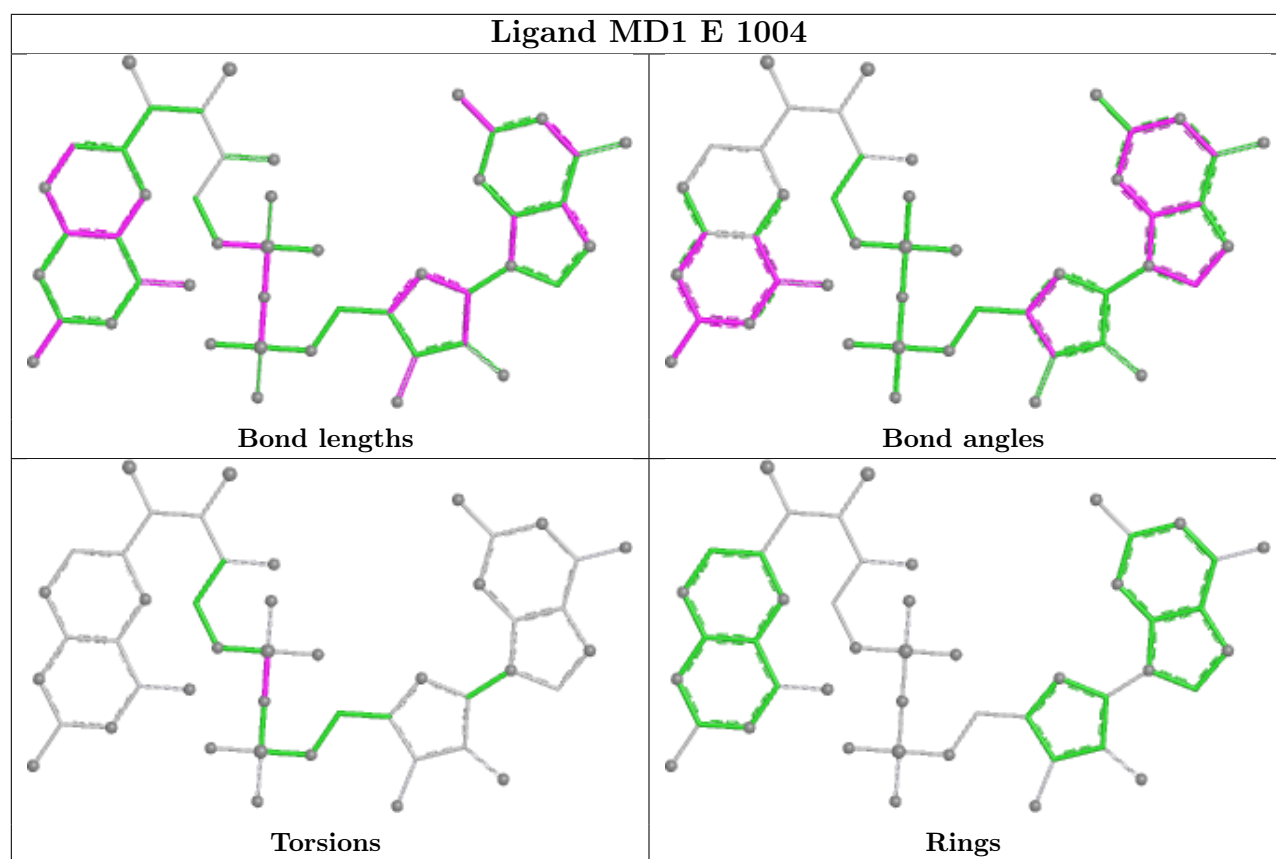
Mol	Chain	Res	Type	Atoms
6	C	1004	MD1	C10-O3A-PB-O1B
10	C	1006	GOL	C1-C2-C3-O3
11	C	1008	PEG	O2-C3-C4-O4
7	E	1008	EDO	O1-C1-C2-O2
11	C	1008	PEG	O1-C1-C2-O2

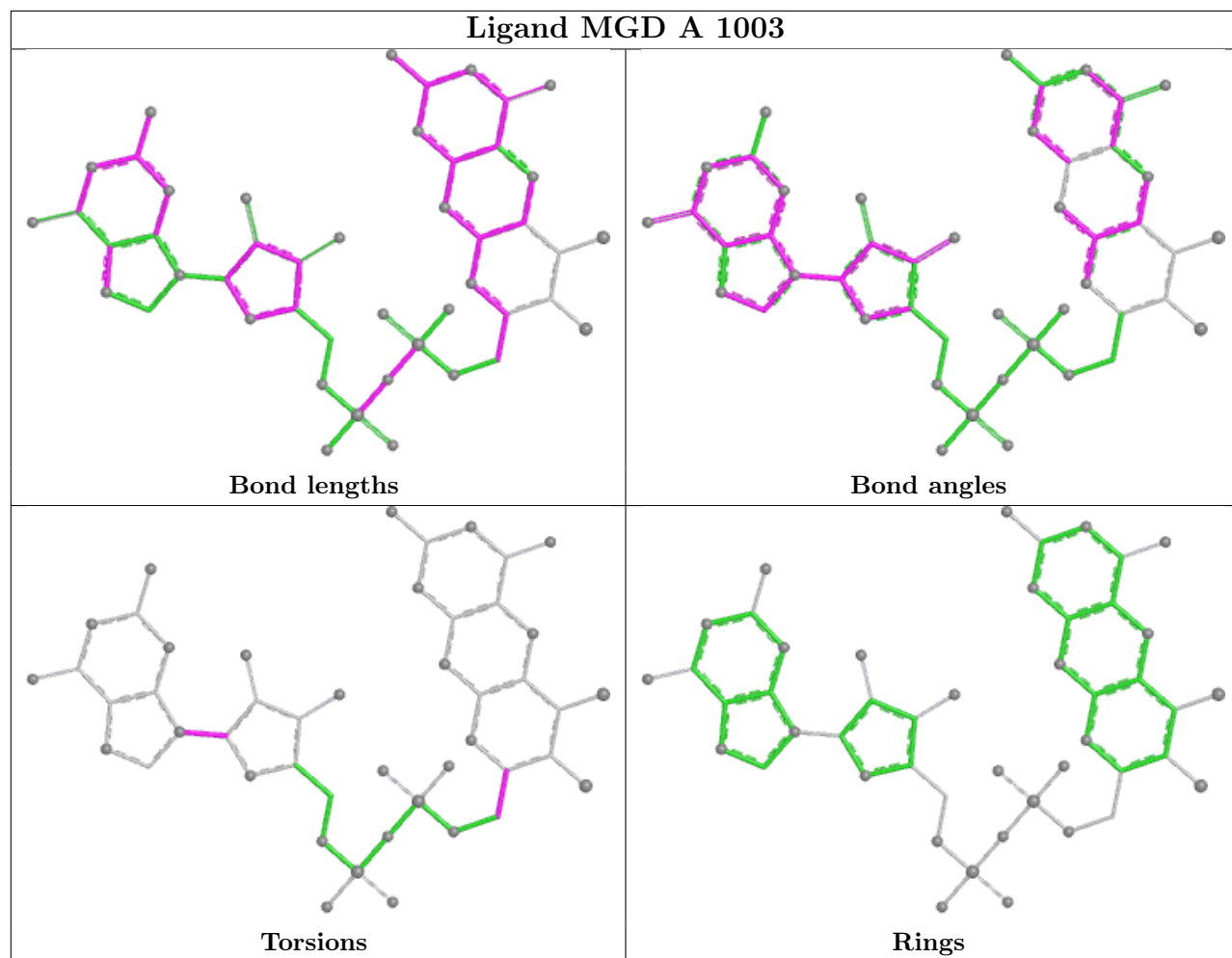
There are no ring outliers.

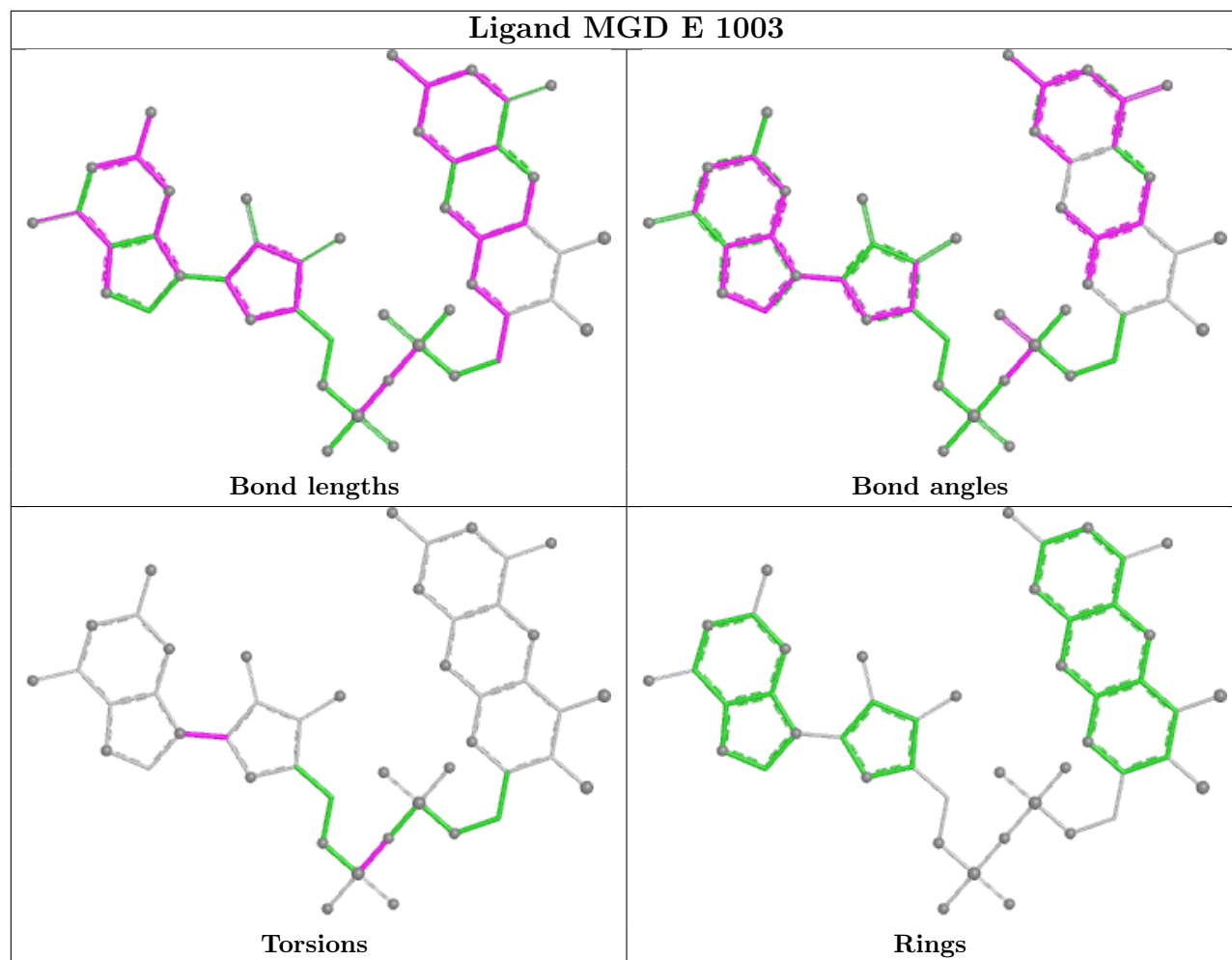
9 monomers are involved in 19 short contacts:

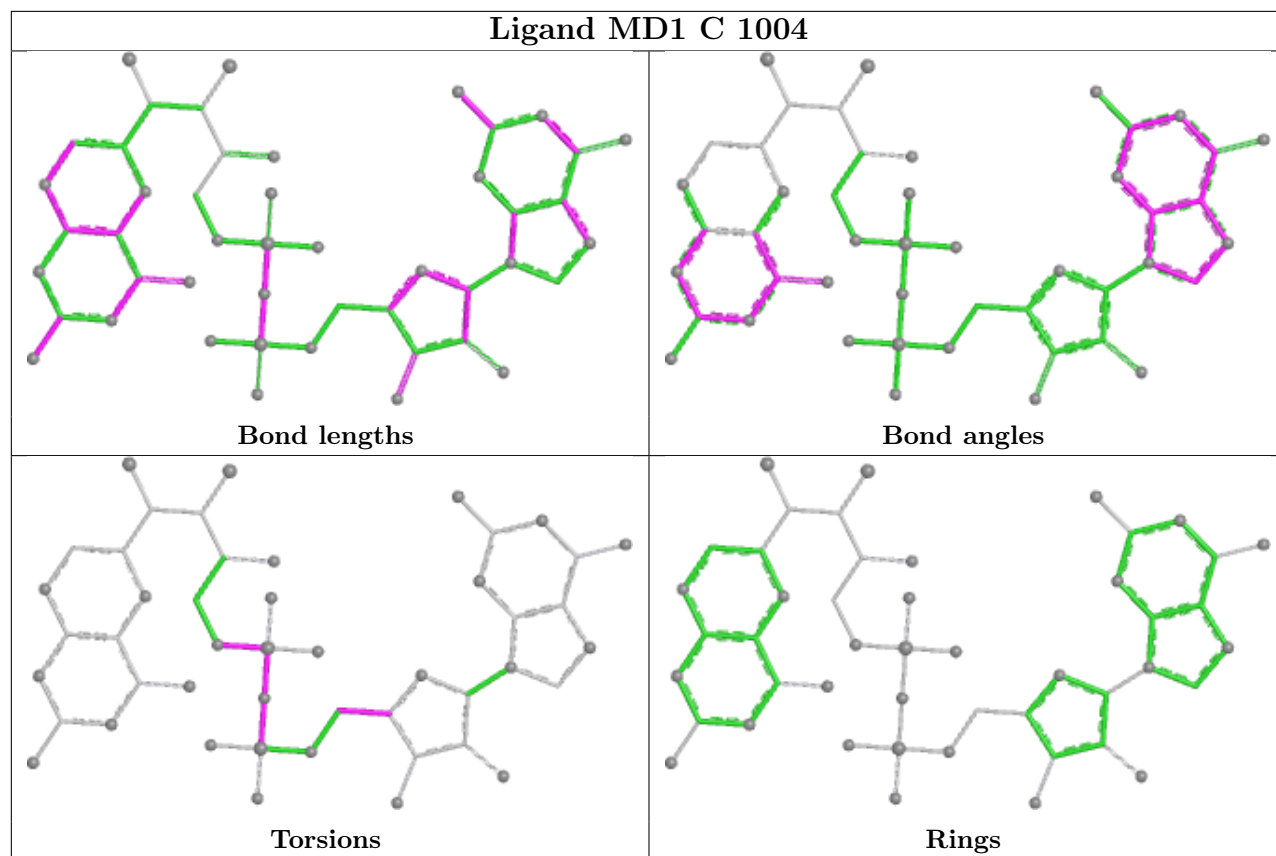
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	E	1004	MD1	4	0
7	A	1010	EDO	1	0
5	E	1003	MGD	1	0
6	C	1004	MD1	4	0
11	C	1008	PEG	1	0
5	C	1003	MGD	1	0
6	A	1004	MD1	5	0
3	D	402	SF4	1	0
8	D	401	F3S	1	0

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.

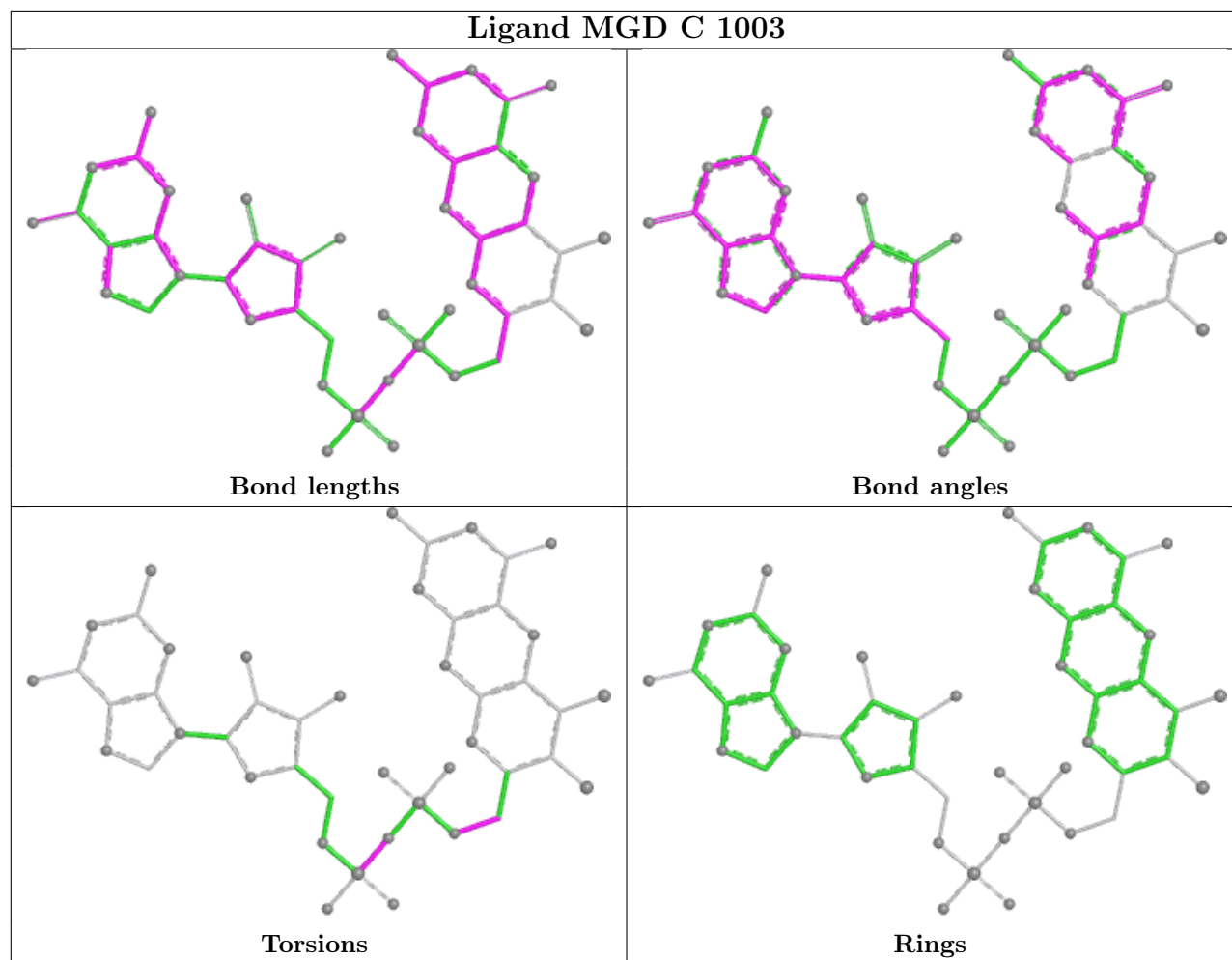


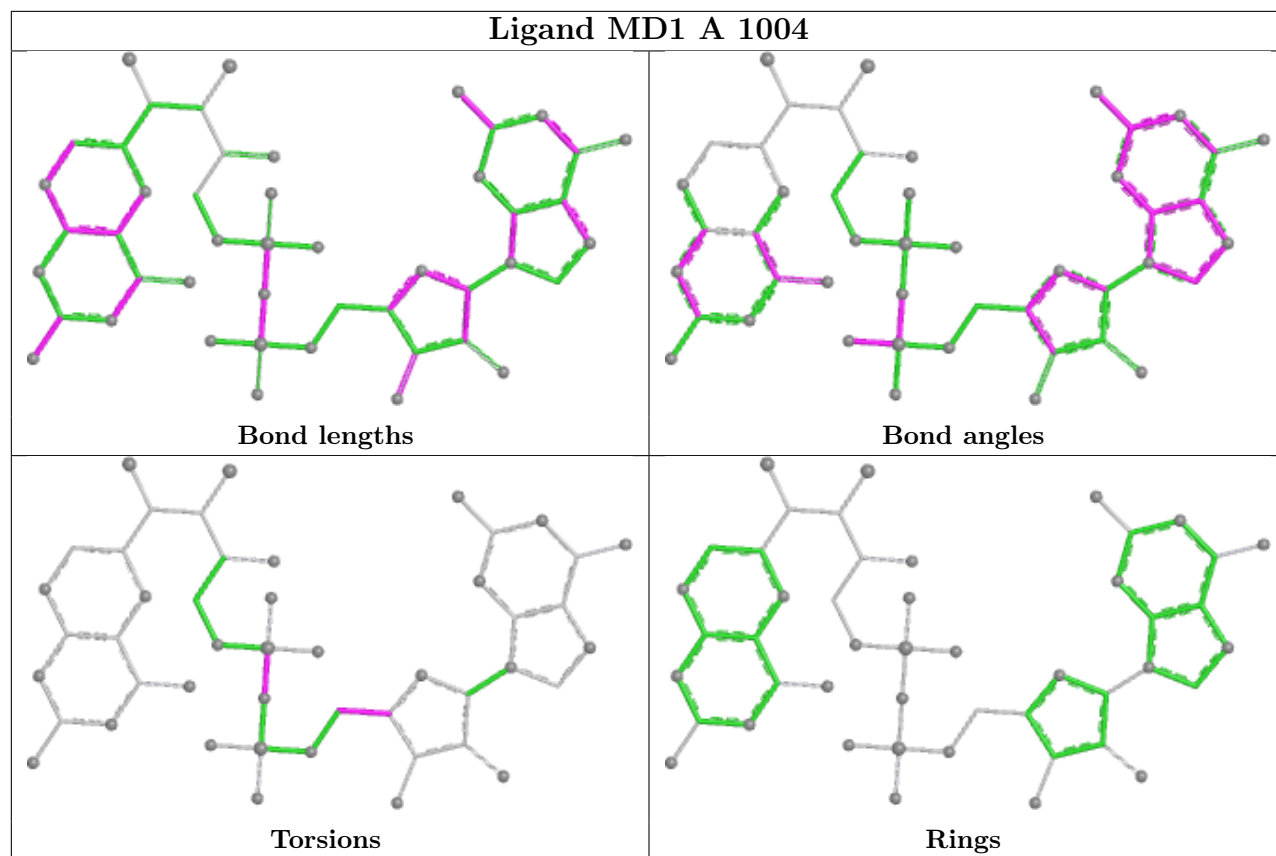






Ligand MGD C 1003





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	895/899 (99%)	0.02	13 (1%)	72 76	11, 26, 46, 73	3 (0%)
1	C	892/899 (99%)	0.48	50 (5%)	30 33	16, 33, 56, 92	0
1	E	892/899 (99%)	0.37	85 (9%)	14 14	14, 27, 60, 84	2 (0%)
2	B	329/333 (98%)	-0.24	2 (0%)	85 88	15, 22, 36, 63	0
2	D	328/333 (98%)	1.69	130 (39%)	1 0	23, 50, 68, 94	0
2	F	328/333 (98%)	0.27	3 (0%)	81 84	15, 29, 47, 67	1 (0%)
All	All	3664/3696 (99%)	0.37	283 (7%)	19 20	11, 29, 58, 94	6 (0%)

The worst 5 of 283 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	282	VAL	6.7
1	A	5	ILE	6.0
1	E	11[A]	TYR	5.9
1	E	360	PHE	5.6
2	D	145	CYS	5.3

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 ⓘ

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
10	GOL	C	1006	6/6	0.69	0.18	37,47,56,57	0
7	EDO	E	1006	4/4	0.72	0.20	25,30,34,36	0
7	EDO	E	1007	4/4	0.74	0.18	25,30,35,35	0
7	EDO	C	1005	4/4	0.80	0.15	27,33,38,42	0
7	EDO	B	406	4/4	0.80	0.17	32,39,40,42	0
7	EDO	E	1008	4/4	0.81	0.15	26,31,35,41	0
7	EDO	E	1009	4/4	0.82	0.18	30,36,40,42	0
11	PEG	C	1008	7/7	0.82	0.15	32,39,42,42	0
7	EDO	A	1010	4/4	0.83	0.20	21,27,30,33	0
7	EDO	A	1008	4/4	0.84	0.14	30,36,41,42	0
7	EDO	E	1010	4/4	0.84	0.14	34,41,43,45	0
7	EDO	A	1009	4/4	0.86	0.13	29,35,41,50	0
9	NA	B	407	1/1	0.87	0.11	35,35,35,35	0
7	EDO	E	1005	4/4	0.90	0.11	16,20,23,23	0
3	SF4	D	404	8/8	0.91	0.09	29,38,39,42	0
7	EDO	C	1007	4/4	0.92	0.09	17,22,26,27	0
7	EDO	A	1007	4/4	0.92	0.09	22,27,28,33	0
3	SF4	D	402	8/8	0.93	0.08	32,42,49,51	0
7	EDO	A	1005	4/4	0.94	0.09	18,25,32,32	0
7	EDO	D	405	4/4	0.94	0.07	26,31,35,35	0
8	F3S	D	401	7/7	0.94	0.08	45,47,55,57	0
5	MGD	E	1003	47/47	0.95	0.08	15,21,29,31	0
3	SF4	F	402	8/8	0.95	0.06	19,22,23,25	0
7	EDO	A	1006	4/4	0.95	0.10	24,29,32,32	0
7	EDO	B	405	4/4	0.96	0.06	16,19,19,20	0
5	MGD	C	1003	47/47	0.96	0.07	20,22,27,30	0
3	SF4	F	404	8/8	0.96	0.06	17,18,20,23	0
5	MGD	A	1003	47/47	0.97	0.06	13,16,19,23	0
7	EDO	F	405	4/4	0.97	0.05	16,19,21,21	0
6	MD1	A	1004	47/47	0.97	0.05	15,17,22,23	0
8	F3S	F	401	7/7	0.97	0.04	21,22,23,24	0
6	MD1	C	1004	47/47	0.97	0.06	17,20,24,26	0
6	MD1	E	1004	47/47	0.97	0.05	12,16,21,23	0
3	SF4	D	403	8/8	0.97	0.06	30,35,39,40	0
3	SF4	B	403	8/8	0.98	0.04	12,14,15,17	0
3	SF4	C	1001	8/8	0.98	0.05	24,25,27,27	0
3	SF4	E	1001	8/8	0.98	0.04	13,15,15,15	0

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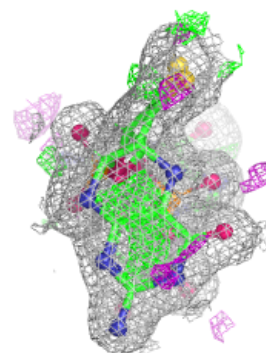
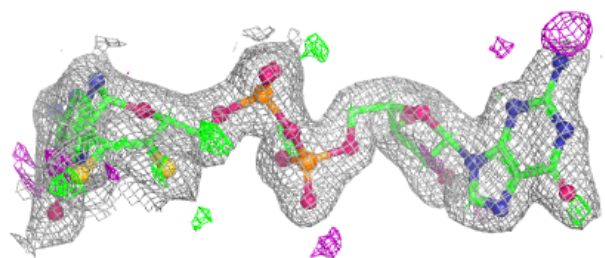
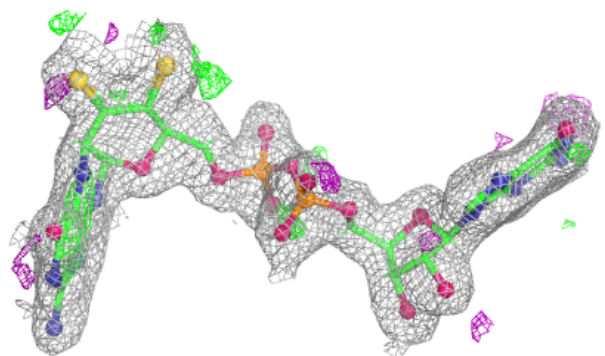
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SF4	B	402	8/8	0.98	0.04	15,16,19,20	0
3	SF4	F	403	8/8	0.98	0.05	18,21,24,25	0
8	F3S	B	401	7/7	0.99	0.03	19,20,22,22	0
4	MO	E	1002	1/1	0.99	0.02	21,21,21,21	0
3	SF4	A	1001	8/8	0.99	0.04	13,14,17,19	0
3	SF4	B	404	8/8	0.99	0.03	12,13,14,15	0
4	MO	A	1002	1/1	0.99	0.03	20,20,20,20	0
4	MO	C	1002	1/1	0.99	0.02	24,24,24,24	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.

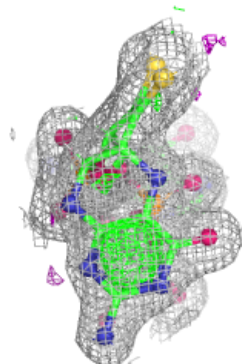
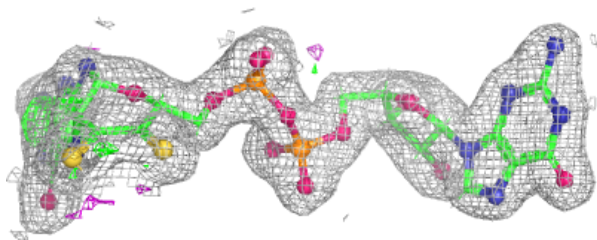
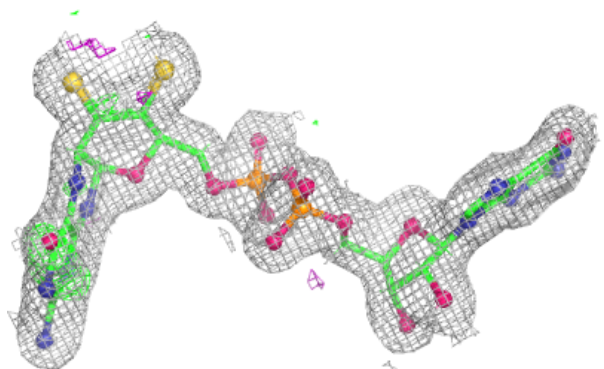
Electron density around MGD E 1003:

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

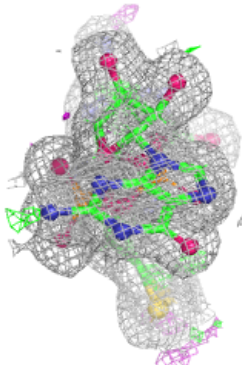
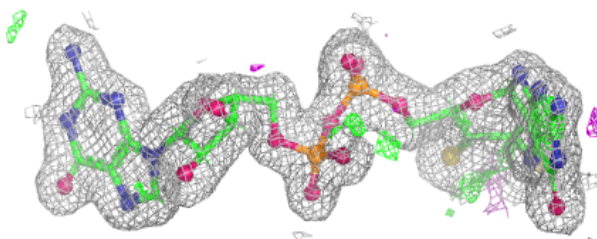
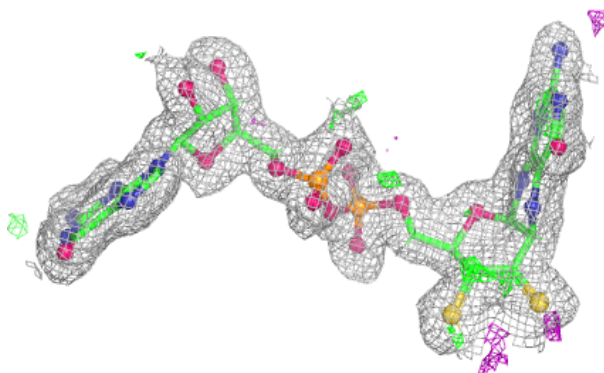


Electron density around MGD C 1003:

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

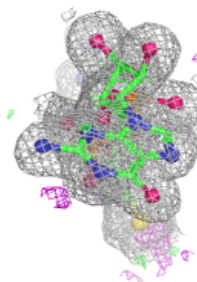
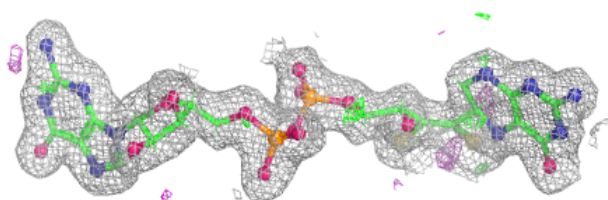
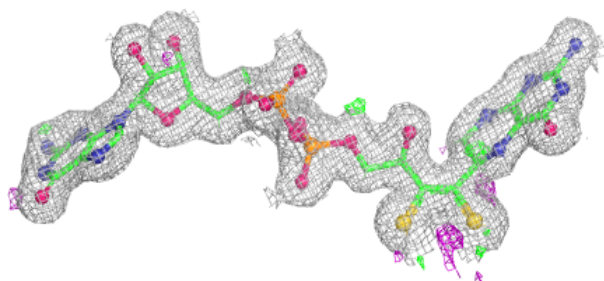
**Electron density around MGD A 1003:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

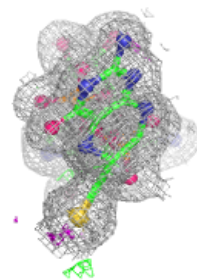
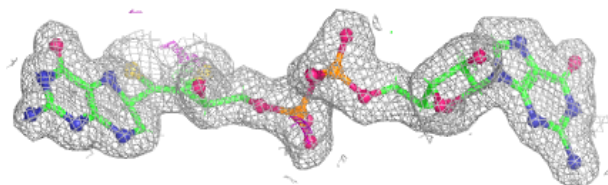
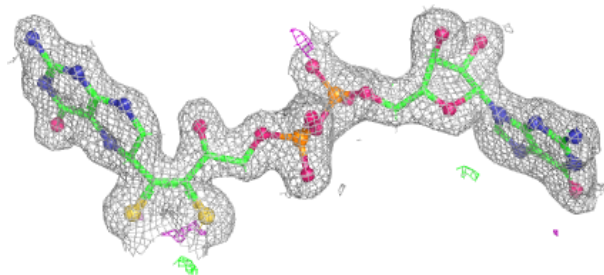


Electron density around MD1 A 1004:

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

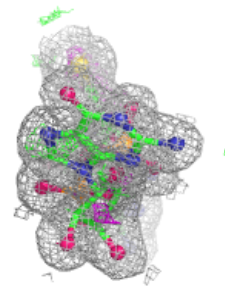
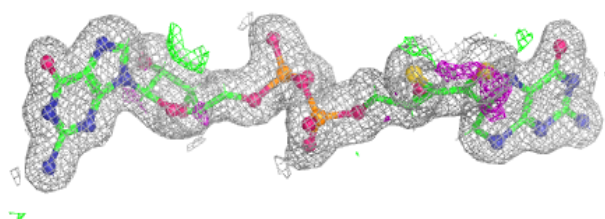
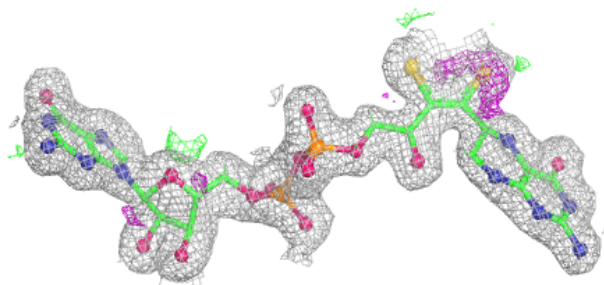
**Electron density around MD1 C 1004:**

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



Electron density around MD1 E 1004:

$2mF_o - DF_c$ (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.