



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:30 PM UTC

PDB ID : 7YZQ / pdb_00007yzq
Title : MgADP-AlF₄-bound DCCP:DCCP-R complex
Authors : Jeoung, J.-H.; Dobbek, H.
Deposited on : 2022-02-21
Resolution : 1.96 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

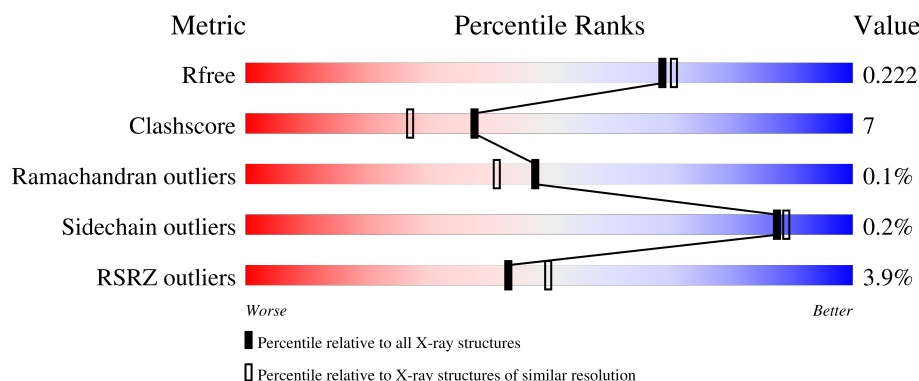
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.96 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.



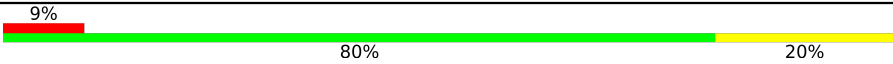
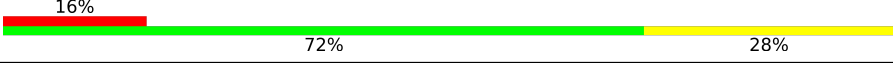
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	422	
1	B	422	
1	C	422	
1	D	422	
2	E	243	

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Mol	Chain	Length	Quality of chain
2	F	243	
2	G	243	
2	H	243	

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 46029 atoms, of which 21708 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 Dehydratase family protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	C	420	Total	C	H	N	O	S	0	3	0
			6776	2157	3402	579	621	17			
1	D	421	Total	C	H	N	O	S	0	6	0
			6886	2184	3471	592	621	18			
1	A	421	Total	C	H	N	O	S	0	6	0
			6828	2172	3429	582	626	19			
1	B	422	Total	C	H	N	O	S	0	10	0
			6900	2189	3472	593	628	18			

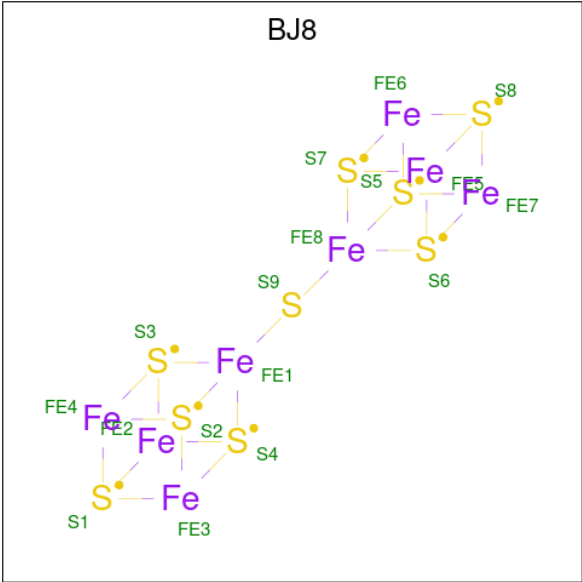
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	GLY	-	expression tag	UNP Q3AET9
D	0	GLY	-	expression tag	UNP Q3AET9
A	0	GLY	-	expression tag	UNP Q3AET9
B	0	GLY	-	expression tag	UNP Q3AET9

- Molecule 2 is a protein called Putative CoA-substrate-specific enzyme activase.

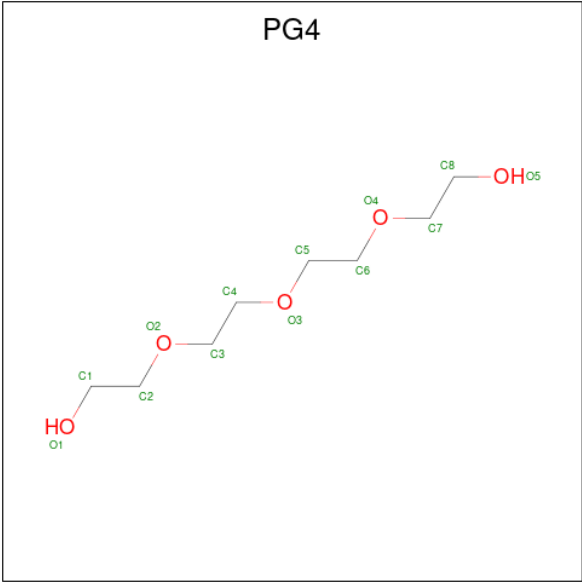
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	E	243	Total	C	H	N	O	S	0	4	0
			3823	1193	1955	321	346	8			
2	F	243	Total	C	H	N	O	S	0	2	0
			3803	1187	1943	320	345	8			
2	G	243	Total	C	H	N	O	S	0	0	0
			3765	1177	1917	318	345	8			
2	H	243	Total	C	H	N	O	S	0	1	0
			3777	1181	1923	318	347	8			

- Molecule 3 is Double cubane cluster (CCD ID: BJ8) (formula: Fe₈S₉) (labeled as "Ligand of Interest" by depositor).



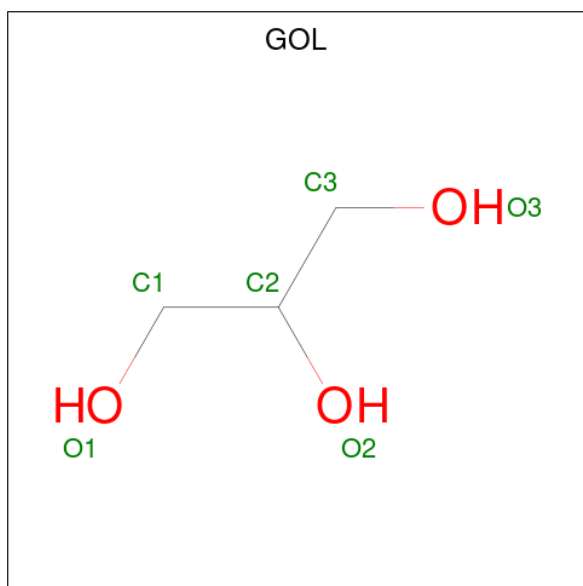
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	Fe	S	0	0
			17	8	9		
3	D	1	Total	Fe	S	0	0
			17	8	9		
3	A	1	Total	Fe	S	0	0
			17	8	9		
3	B	1	Total	Fe	S	0	0
			17	8	9		

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



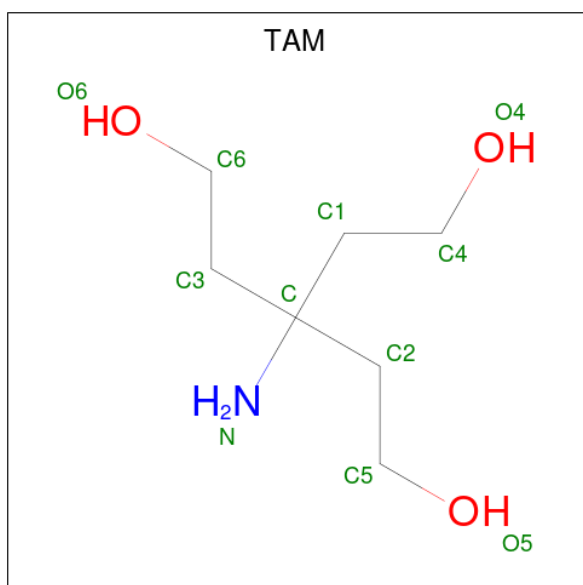
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	H	O	0	1
			62	16	36	10		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



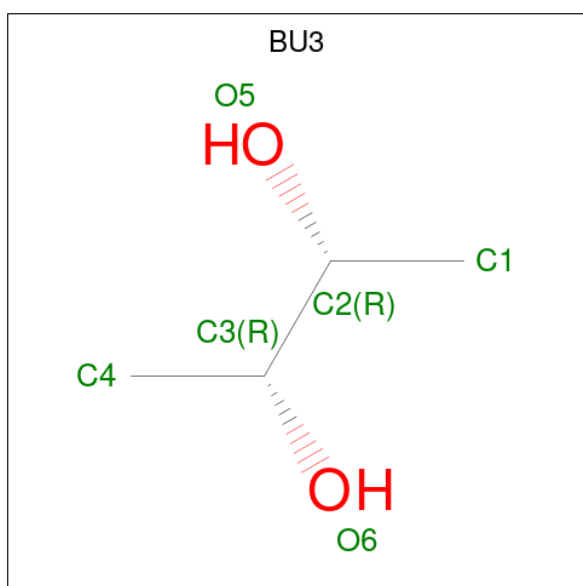
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 6 is TRIS(HYDROXYETHYL)AMINOMETHANE (CCD ID: TAM) (formula: $C_7H_{17}NO_3$).



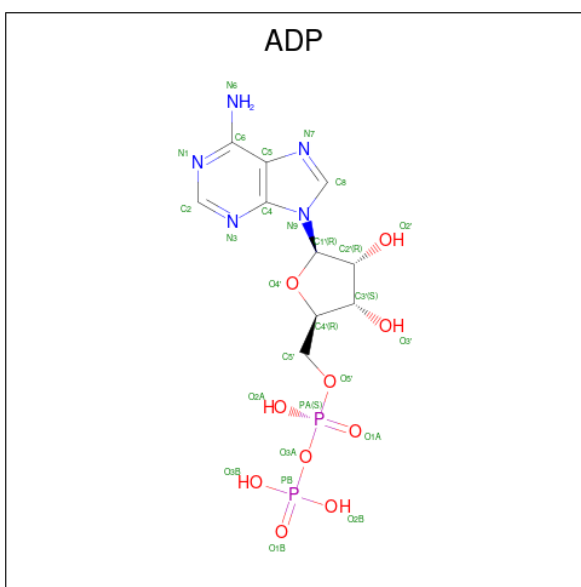
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	D	1	Total 28	C 7	H 17	N 1	O 3	0	0
6	E	1	Total 28	C 7	H 17	N 1	O 3	0	0
6	B	1	Total 28	C 7	H 17	N 1	O 3	0	0
6	G	1	Total 28	C 7	H 17	N 1	O 3	0	0

- Molecule 7 is (R,R)-2,3-BUTANEDIOL (CCD ID: BU3) (formula: $C_4H_{10}O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	D	1	Total	C	H	O	0	0
			16	4	10	2		
7	B	1	Total	C	H	O	0	0
			16	4	10	2		
7	B	1	Total	C	H	O	0	0
			16	4	10	2		
7	G	1	Total	C	H	O	0	0
			16	4	10	2		

- Molecule 8 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

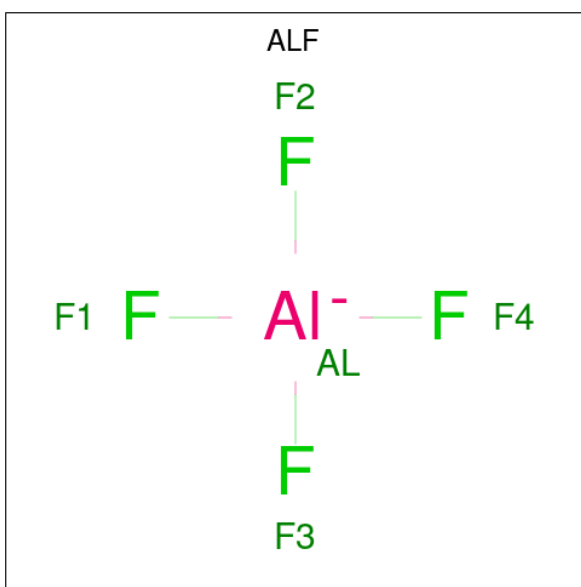


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	E	1	Total	C	H	N	O	P	
			38	10	11	5	10	2	0
8	F	1	Total	C	H	N	O	P	
			38	10	11	5	10	2	0
8	G	1	Total	C	H	N	O	P	
			38	10	11	5	10	2	0
8	H	1	Total	C	H	N	O	P	
			38	10	11	5	10	2	0

- Molecule 9 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

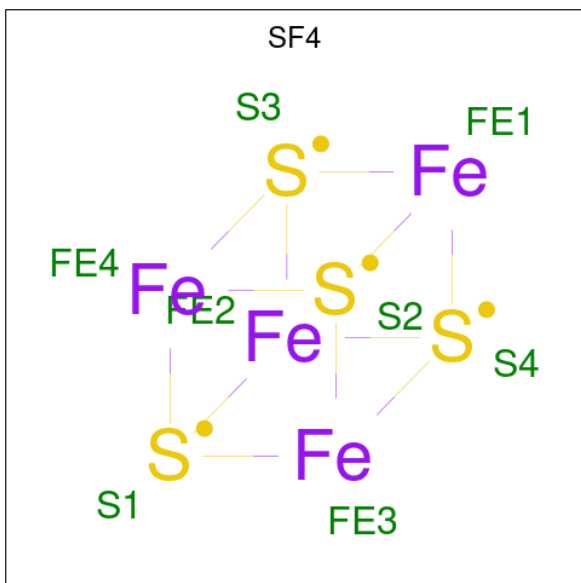
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	E	1	Total	Mg		
			1	1	0	0
9	F	1	Total	Mg		
			1	1	0	0
9	G	1	Total	Mg		
			1	1	0	0
9	H	1	Total	Mg		
			1	1	0	0

- Molecule 10 is TETRAFLUOROALUMINATE ION (CCD ID: ALF) (formula: AlF_4).



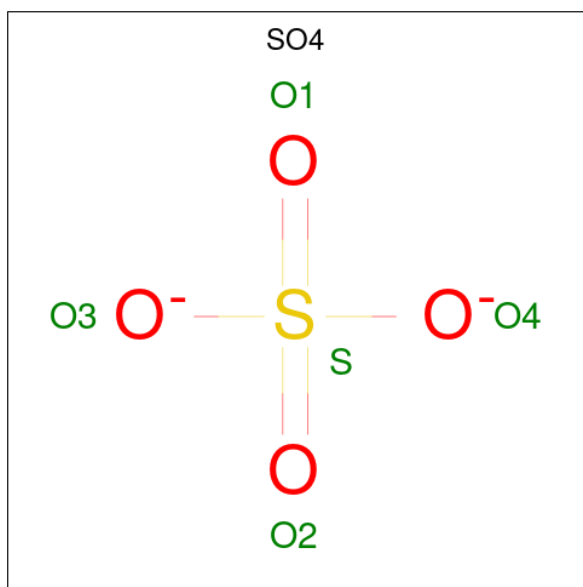
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	E	1	Total	Al	F	0	0
			5	1	4		
10	F	1	Total	Al	F	0	0
			5	1	4		
10	G	1	Total	Al	F	0	0
			5	1	4		
10	H	1	Total	Al	F	0	0
			5	1	4		

- Molecule 11 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



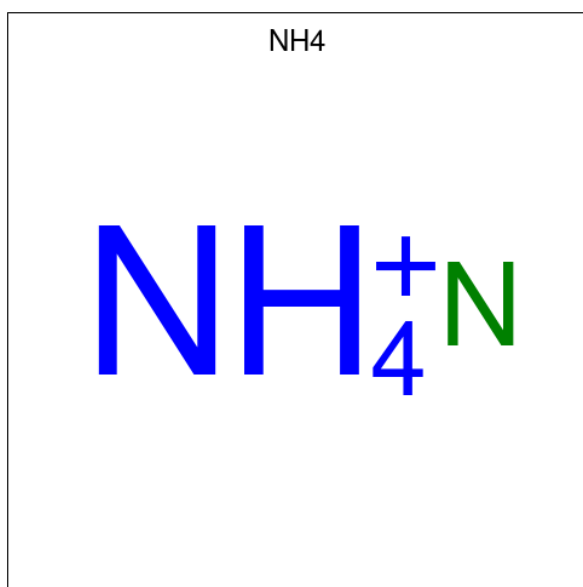
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	F	1	Total	Fe	S	0	0
			8	4	4		
11	G	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 12 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	F	1	Total	O	S	0	0
			5	4	1		
12	A	1	Total	O	S	0	0
			5	4	1		
12	A	1	Total	O	S	0	0
			5	4	1		
12	A	1	Total	O	S	0	0
			5	4	1		
12	G	1	Total	O	S	0	0
			5	4	1		

- Molecule 13 is AMMONIUM ION (CCD ID: NH4) (formula: H₄N).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
13	A	1	Total N 1 1	0	0
13	G	1	Total N 1 1	0	0

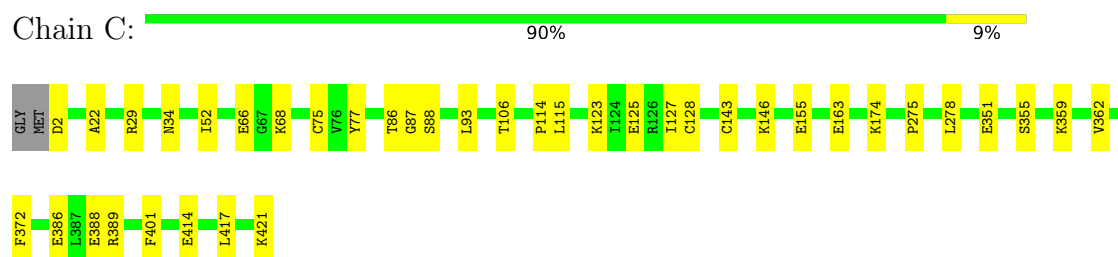
- Molecule 14 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
14	C	584	Total O 584 584	0	0
14	D	435	Total O 435 435	0	0
14	E	256	Total O 256 256	0	0
14	F	154	Total O 154 154	0	0
14	A	632	Total O 632 632	0	0
14	B	441	Total O 441 441	0	0
14	G	271	Total O 271 271	0	0
14	H	159	Total O 159 159	0	0

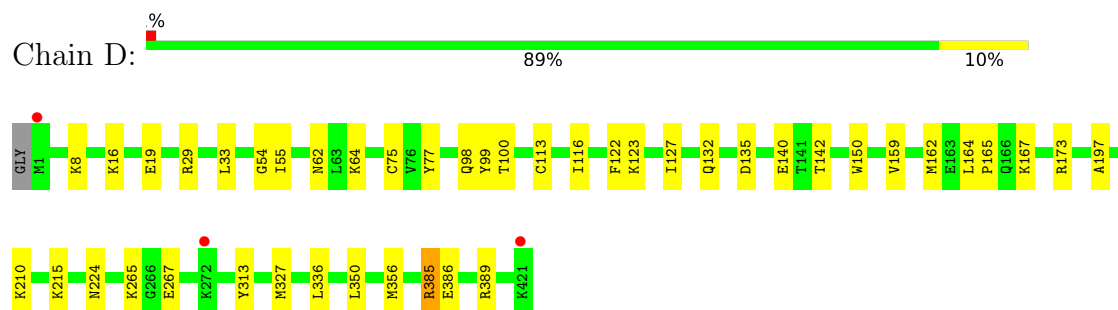
3 Residue-property plots [i](#)

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.

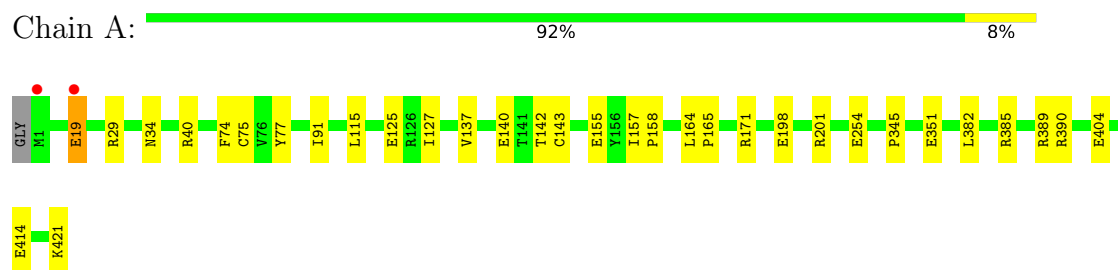
• Molecule 1: Dehydratase family protein



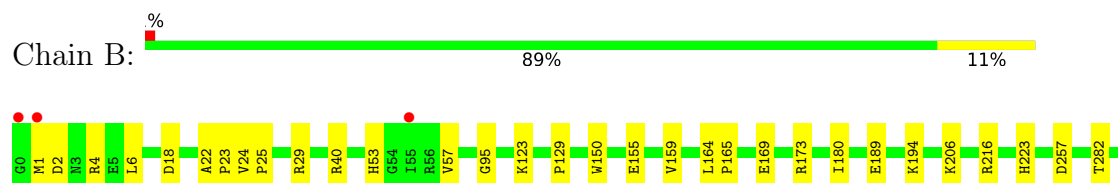
• Molecule 1: Dehydratase family protein



• Molecule 1: Dehydratase family protein

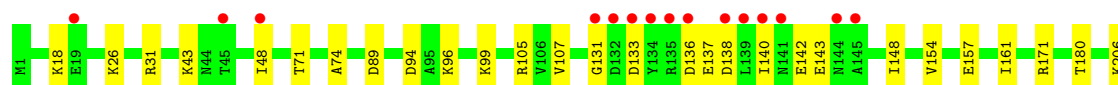
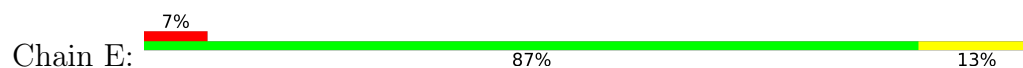


• Molecule 1: Dehydratase family protein

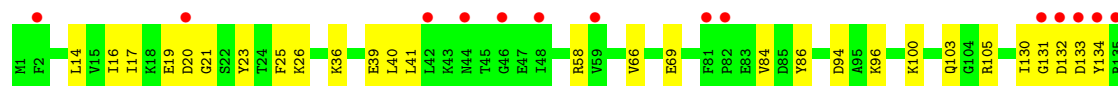
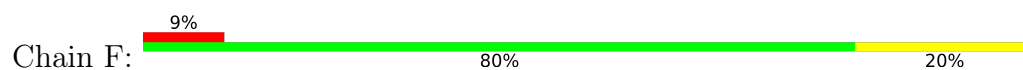




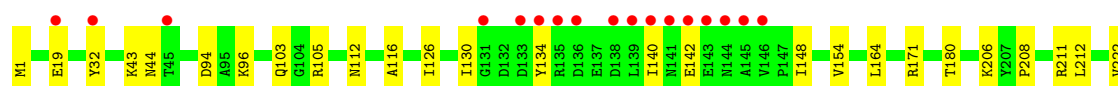
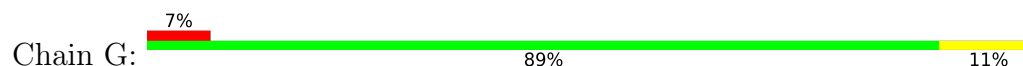
- Molecule 2: Putative CoA-substrate-specific enzyme activase



- Molecule 2: Putative CoA-substrate-specific enzyme activase



- Molecule 2: Putative CoA-substrate-specific enzyme activase



- Molecule 2: Putative CoA-substrate-specific enzyme activase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	78.03Å 81.59Å 121.48Å 100.83° 96.92° 90.17°	Depositor
Resolution (Å)	46.16 – 1.96 46.16 – 1.96	Depositor EDS
% Data completeness (in resolution range)	96.0 (46.16-1.96) 96.4 (46.16-1.96)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.178 , 0.223 0.178 , 0.222	Depositor DCC
R_{free} test set	2128 reflections (1.05%)	wwPDB-VP
Wilson B-factor (Å ²)	19.6	Xtriage
Anisotropy	0.379	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	46029	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.07% 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: ALF, SF4, NH4, SO4, ADP, TAM, GOL, MG, PG4, BU3, BJ8

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	0.68	2/3486 (0.1%)	0.66	0/4707
1	B	0.51	0/3532	0.56	0/4766
1	C	0.65	0/3452	0.69	0/4663
1	D	0.52	2/3505 (0.1%)	0.58	0/4728
2	E	0.70	1/1906 (0.1%)	0.70	1/2572 (0.0%)
2	F	0.45	0/1892	0.57	0/2551
2	G	0.58	0/1874	0.63	0/2528
2	H	0.52	0/1883	0.65	2/2540 (0.1%)
All	All	0.59	5/21530 (0.0%)	0.63	3/29055 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	225	PRO	CA-C	14.27	1.59	1.51
1	A	143	CYS	CB-SG	-9.01	1.51	1.81
1	D	356	MET	SD-CE	-6.29	1.63	1.79
1	A	91	ILE	C-N	6.13	1.36	1.33
1	D	162	MET	SD-CE	-5.12	1.66	1.79

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	225	PRO	O-C-N	10.00	125.91	121.31
2	H	134	TYR	CA-C-N	-6.12	111.81	122.64
2	H	134	TYR	C-N-CA	-6.12	111.81	122.64

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	3399	3429	3431	26	0
1	B	3428	3472	3453	38	0
1	C	3374	3402	3402	40	1
1	D	3415	3471	3473	33	0
2	E	1868	1955	1955	32	0
2	F	1860	1943	1943	47	0
2	G	1848	1917	1917	24	0
2	H	1854	1923	1923	71	0
3	A	17	0	0	0	0
3	B	17	0	0	0	0
3	C	17	0	0	1	0
3	D	17	0	0	0	0
4	C	26	36	36	5	0
5	D	6	8	8	0	0
6	B	11	17	17	1	0
6	D	11	17	17	0	0
6	E	11	17	17	0	0
6	G	11	17	17	0	0
7	B	12	20	20	3	0
7	D	6	10	10	0	0
7	G	6	10	10	0	0
8	E	27	11	12	0	0
8	F	27	11	12	0	0
8	G	27	11	12	0	0
8	H	27	11	12	1	0
9	E	1	0	0	0	0
9	F	1	0	0	0	0
9	G	1	0	0	0	0
9	H	1	0	0	0	0
10	E	5	0	0	0	0
10	F	5	0	0	0	0
10	G	5	0	0	0	0
10	H	5	0	0	0	0
11	F	8	0	0	0	0
11	G	8	0	0	0	0
12	A	15	0	0	1	0
12	F	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	G	5	0	0	1	0
13	A	1	0	0	1	0
13	G	1	0	0	1	0
14	A	632	0	0	15	4
14	B	441	0	0	9	0
14	C	584	0	0	17	3
14	D	435	0	0	4	0
14	E	256	0	0	12	2
14	F	154	0	0	5	0
14	G	271	0	0	5	4
14	H	159	0	0	11	0
All	All	24321	21708	21697	303	7

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (303) 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)
2:H:138:ASP:HB3	2:H:171:ARG:NH1	1.64	1.12
2:H:179:LYS:NZ	14:H:401:HOH:O	1.90	1.04
2:F:206:LYS:NZ	14:F:402:HOH:O	1.84	1.01
2:G:105:ARG:NH2	14:G:401:HOH:O	1.93	1.00
1:B:4:ARG:NH2	1:B:18:ASP:OD2	1.99	0.94
2:H:140:ILE:HD13	2:H:213:PHE:CE1	2.04	0.93
2:H:140:ILE:HD13	2:H:213:PHE:HE1	1.31	0.93
2:F:36:LYS:NZ	14:F:401:HOH:O	1.84	0.90
2:H:138:ASP:O	2:H:171:ARG:NH1	2.07	0.88
1:C:351:GLU:OE1	14:C:601:HOH:O	1.90	0.87
1:C:417:LEU:O	1:C:421:LYS:HD3	1.75	0.87
2:E:26:LYS:NZ	14:E:401:HOH:O	2.07	0.87
1:C:88:SER:OG	14:C:602:HOH:O	1.92	0.86
2:E:105:ARG:NH2	14:E:403:HOH:O	2.08	0.84
1:D:164:LEU:HD12	1:D:165:PRO:HD2	1.60	0.84
1:C:155:GLU:OE2	1:D:385[A]:ARG:NH2	2.11	0.82
2:G:103:GLN:O	14:G:402:HOH:O	1.96	0.82
1:D:350:LEU:HD21	1:D:386:GLU:HG2	1.62	0.81
2:H:138:ASP:CB	2:H:171:ARG:NH1	2.45	0.80
2:G:171:ARG:NH1	14:G:403:HOH:O	2.15	0.79
1:C:414:GLU:OE1	14:C:603:HOH:O	2.00	0.78
2:G:134:TYR:HE1	2:G:171:ARG:HB2	1.47	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:142:GLU:OE1	2:H:171:ARG:NE	2.17	0.78
2:G:32:TYR:HB3	2:H:164:LEU:HD23	1.65	0.78
2:H:140:ILE:CD1	2:H:213:PHE:HE1	1.96	0.77
1:D:123[C]:LYS:NZ	1:D:132:GLN:OE1	2.17	0.77
1:A:19[B]:GLU:OE2	14:A:601:HOH:O	2.02	0.76
12:A:501:SO4:O1	13:A:502:NH4:N	2.18	0.76
2:H:138:ASP:C	2:H:171:ARG:NH1	2.42	0.76
1:B:206:LYS:NZ	14:B:603:HOH:O	2.20	0.75
12:G:304:SO4:O4	13:G:305:NH4:N	2.20	0.75
2:F:138:ASP:HB3	2:F:171:ARG:HE	1.52	0.74
2:H:138:ASP:HB3	2:H:171:ARG:HH12	1.50	0.74
2:H:84:VAL:O	2:H:100:LYS:NZ	2.21	0.72
2:G:211:ARG:NH2	2:G:222:VAL:O	2.21	0.72
1:B:189:GLU:OE2	14:B:601:HOH:O	2.09	0.70
1:A:351:GLU:OE2	14:A:603:HOH:O	2.09	0.70
1:A:414:GLU:OE1	14:A:602:HOH:O	2.08	0.70
2:H:31:ARG:NH2	14:H:405:HOH:O	2.24	0.70
2:H:138:ASP:CB	2:H:171:ARG:HH12	2.04	0.69
1:A:254:GLU:OE1	14:A:604:HOH:O	2.10	0.69
2:H:135:ARG:NH2	14:H:402:HOH:O	2.06	0.69
2:H:36:LYS:H	2:H:36:LYS:HD2	1.57	0.68
2:F:40:LEU:N	14:F:401:HOH:O	2.25	0.68
2:G:44:ASN:ND2	14:G:404:HOH:O	2.27	0.67
1:C:388:GLU:OE2	14:C:604:HOH:O	2.11	0.67
2:H:36:LYS:HD2	2:H:36:LYS:N	2.09	0.67
1:C:22:ALA:O	14:C:605:HOH:O	2.13	0.67
1:D:385[A]:ARG:NH1	14:D:602:HOH:O	2.28	0.66
2:E:105:ARG:NH1	14:E:406:HOH:O	2.29	0.66
2:E:74:ALA:HB3	14:E:403:HOH:O	1.95	0.65
2:H:85:ASP:OD1	14:H:403:HOH:O	2.14	0.65
2:H:170:LYS:NZ	14:H:409:HOH:O	2.31	0.64
1:A:171:ARG:NE	14:A:614:HOH:O	2.30	0.63
1:D:385[A]:ARG:HH11	1:D:385[A]:ARG:HB3	1.63	0.63
1:B:223:HIS:HA	7:B:503:BU3:H3	1.79	0.63
2:H:138:ASP:C	2:H:171:ARG:HH11	2.07	0.63
1:B:40[B]:ARG:NH1	14:B:605:HOH:O	2.27	0.62
2:H:138:ASP:CA	2:H:171:ARG:HH12	2.12	0.62
2:H:206:LYS:HE3	2:H:229:SER:OG	1.99	0.62
2:E:71:THR:HG23	2:E:105:ARG:HH12	1.63	0.61
1:B:392:ILE:O	14:B:602:HOH:O	2.16	0.61
2:F:14:LEU:HD11	2:F:16:ILE:HD11	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:33:GLU:HB3	2:H:36:LYS:HD3	1.82	0.61
2:H:99:LYS:NZ	14:H:410:HOH:O	2.34	0.60
1:A:40:ARG:NH1	14:A:616:HOH:O	2.33	0.60
1:A:421:LYS:NZ	14:A:618:HOH:O	2.34	0.59
2:H:136:ASP:OD1	2:H:209:ALA:HB2	2.02	0.59
2:E:142:GLU:HG2	2:E:171:ARG:HD3	1.84	0.59
2:F:140:ILE:N	2:F:140:ILE:HD12	2.18	0.59
2:F:148:ILE:O	2:F:180:THR:HG21	2.02	0.59
2:G:134:TYR:HE1	2:G:171:ARG:CB	2.16	0.59
1:D:123[A]:LYS:HA	1:D:123[A]:LYS:HE3	1.85	0.58
2:H:166:ARG:NH1	14:H:407:HOH:O	2.27	0.58
2:H:187:LYS:O	2:H:190[B]:GLU:HG3	2.03	0.58
2:F:140:ILE:HG12	2:F:212:LEU:HD12	1.85	0.58
1:B:6:LEU:HD21	1:B:57:VAL:HG12	1.84	0.57
1:D:62:ASN:OD1	14:D:601:HOH:O	2.17	0.57
2:H:17:ILE:HG13	2:H:237:ALA:HB1	1.86	0.57
2:E:94:ASP:HB2	2:E:96:LYS:HE3	1.86	0.57
1:B:150:TRP:HB3	1:B:159:VAL:HG11	1.87	0.57
2:H:36:LYS:H	2:H:36:LYS:CD	2.17	0.57
1:C:421:LYS:NZ	14:C:612:HOH:O	2.30	0.56
2:E:138:ASP:OD2	2:E:171:ARG:NH2	2.38	0.56
1:A:198:GLU:OE2	14:A:605:HOH:O	2.18	0.56
2:G:1:MET:HE1	2:G:19:GLU:HB2	1.86	0.56
2:G:154:VAL:HG23	2:H:154:VAL:HG23	1.87	0.56
2:F:86:TYR:CD1	2:F:195:PRO:HB3	2.42	0.55
2:H:148:ILE:O	2:H:180:THR:HG21	2.07	0.55
1:C:93:LEU:HB3	4:C:502[A]:PG4:H71	1.89	0.55
2:E:154:VAL:HG23	2:F:154:VAL:HG23	1.89	0.55
1:B:180:ILE:HG13	1:B:332:ALA:HB2	1.88	0.55
1:B:385[B]:ARG:O	1:B:389:ARG:HG3	2.07	0.55
1:C:125:GLU:HB2	1:C:127:ILE:HG13	1.89	0.54
2:H:1:MET:N	2:H:17:ILE:O	2.40	0.54
1:C:2:ASP:OD2	14:C:606:HOH:O	2.18	0.54
2:F:21:GLY:HA2	2:F:241:ARG:NH2	2.23	0.54
1:D:16:LYS:HB3	1:D:127:ILE:HD11	1.90	0.54
1:D:197:ALA:HA	1:D:327:MET:HE1	1.90	0.54
1:C:163:GLU:HB2	1:D:167:LYS:HE3	1.90	0.54
1:C:128:CYS:SG	4:C:502[A]:PG4:H61	2.48	0.53
2:E:96:LYS:HD2	2:E:96:LYS:N	2.23	0.53
2:H:142:GLU:OE1	2:H:171:ARG:CD	2.56	0.53
1:D:8:LYS:NZ	14:D:620:HOH:O	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:385[A]:ARG:CZ	1:D:389:ARG:HH12	2.21	0.53
2:F:58:ARG:HD2	2:F:66:VAL:HG13	1.90	0.53
2:H:126:ILE:HG21	2:H:177:LEU:HD13	1.91	0.52
2:H:216:LYS:NZ	14:H:415:HOH:O	2.41	0.52
1:C:68:LYS:HE2	1:C:86:THR:O	2.10	0.52
1:C:123:LYS:NZ	14:C:630:HOH:O	2.43	0.52
1:B:315:LYS:NZ	14:B:615:HOH:O	2.42	0.52
2:F:138:ASP:O	2:F:171:ARG:NH1	2.39	0.52
2:E:105:ARG:CZ	2:E:105:ARG:HA	2.40	0.51
1:A:155:GLU:OE2	1:B:385[B]:ARG:NH1	2.43	0.51
2:G:94:ASP:HB2	2:G:96:LYS:HE3	1.92	0.51
2:H:218:MET:HB3	2:H:220:VAL:HG22	1.91	0.51
1:C:143:CYS:SG	1:C:146:LYS:HB2	2.50	0.51
2:H:179:LYS:HD3	14:H:459:HOH:O	2.11	0.50
14:C:722:HOH:O	2:F:151:MET:CE	2.58	0.50
2:F:26:LYS:HG3	2:F:41:LEU:CD2	2.41	0.50
2:F:194[B]:LYS:HG3	14:F:443:HOH:O	2.09	0.50
1:D:210:LYS:HE3	1:D:313:TYR:OH	2.11	0.50
1:C:278:LEU:HB2	1:C:362:VAL:HG21	1.93	0.50
2:E:99:LYS:O	2:E:107[A]:VAL:HG22	2.11	0.50
2:E:148:ILE:H	2:E:180[B]:THR:HG21	1.76	0.50
1:C:372:PHE:HE2	2:F:151:MET:HE3	1.77	0.50
1:C:275:PRO:HG3	1:C:421:LYS:HE3	1.94	0.49
1:D:265:LYS:HB2	1:D:267:GLU:HG3	1.94	0.49
1:C:421:LYS:NZ	14:C:617:HOH:O	2.35	0.49
1:C:68:LYS:NZ	14:C:602:HOH:O	2.09	0.49
1:D:224:ASN:ND2	14:D:610:HOH:O	2.36	0.49
2:F:134:TYR:HE2	2:F:170:LYS:HB2	1.77	0.49
2:H:130:ILE:HG21	2:H:164:LEU:HD21	1.94	0.49
2:H:138:ASP:CA	2:H:171:ARG:NH1	2.72	0.49
1:C:388:GLU:HB2	14:C:896:HOH:O	2.13	0.49
1:B:223:HIS:ND1	7:B:503:BU3:H3	2.27	0.49
1:C:68:LYS:HD3	14:C:671:HOH:O	2.12	0.49
6:B:504:TAM:H41	14:B:769:HOH:O	2.13	0.49
2:H:78:HIS:HA	2:H:100:LYS:HE2	1.95	0.49
2:H:137:GLU:O	2:H:141:ASN:OD1	2.30	0.49
2:H:164:LEU:N	2:H:164:LEU:HD12	2.28	0.48
1:C:275:PRO:HD3	1:C:421:LYS:HE2	1.95	0.48
2:H:190[B]:GLU:OE1	14:H:404:HOH:O	2.20	0.48
2:F:84:VAL:HG11	2:F:199:ILE:HG13	1.95	0.48
1:B:2:ASP:N	1:B:2:ASP:OD1	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:CYS:HB2	4:C:502[A]:PG4:H61	1.94	0.48
2:F:133:ASP:OD1	2:F:133:ASP:C	2.56	0.47
2:G:96:LYS:HD2	2:G:96:LYS:N	2.29	0.47
2:H:94:ASP:HB2	2:H:96:LYS:HE3	1.96	0.47
1:D:140:GLU:HG2	1:D:142:THR:HG23	1.96	0.47
2:F:138:ASP:HB3	2:F:171:ARG:NE	2.23	0.47
1:A:390:ARG:NH2	14:A:639:HOH:O	2.47	0.47
1:B:194:LYS:NZ	14:B:621:HOH:O	2.47	0.47
2:E:148:ILE:O	2:E:180[A]:THR:HG21	2.15	0.47
2:G:206:LYS:O	2:G:208:PRO:HD3	2.14	0.47
1:C:75:CYS:SG	1:C:77:TYR:HB2	2.54	0.47
2:E:211[B]:ARG:HD2	2:E:224:VAL:HG23	1.96	0.47
2:F:132:ASP:OD1	2:F:134:TYR:HB2	2.15	0.47
2:G:211:ARG:HD2	2:G:224:VAL:HG23	1.97	0.47
2:G:134:TYR:CD1	2:G:171:ARG:HG3	2.49	0.47
2:F:103:GLN:HB2	2:F:105:ARG:HE	1.80	0.47
2:F:14:LEU:O	2:F:25:PHE:HA	2.15	0.47
2:F:17:ILE:HG12	2:F:23:TYR:HB3	1.96	0.47
1:B:419:MET:HG3	2:H:106:VAL:HG23	1.97	0.47
2:F:94:ASP:HB2	2:F:96:LYS:HE3	1.97	0.46
14:C:722:HOH:O	2:F:151:MET:HE3	2.14	0.46
1:C:115:LEU:C	1:C:115:LEU:HD23	2.40	0.46
1:A:385[B]:ARG:CZ	1:A:389:ARG:HH22	2.29	0.46
1:B:385[A]:ARG:O	1:B:389:ARG:HG3	2.16	0.46
2:E:18:LYS:CD	14:E:414:HOH:O	2.62	0.46
2:E:137:GLU:O	2:E:140:ILE:HG22	2.16	0.46
1:A:385[B]:ARG:NH2	1:B:155[B]:GLU:OE2	2.49	0.46
2:G:148:ILE:O	2:G:180:THR:HG21	2.16	0.46
1:D:150:TRP:HB3	1:D:159:VAL:HG11	1.98	0.45
2:E:180[B]:THR:HG23	14:E:461:HOH:O	2.16	0.45
1:D:19:GLU:HG2	1:D:99:TYR:CE1	2.51	0.45
1:D:385[A]:ARG:HH11	1:D:385[A]:ARG:CB	2.27	0.45
1:B:382:LEU:HG	14:B:711:HOH:O	2.15	0.45
1:C:75:CYS:HA	3:C:501:BJ8:S5	2.57	0.45
2:F:130:ILE:CG2	2:F:170:LYS:HB3	2.47	0.45
1:D:173[B]:ARG:HG2	1:D:336:LEU:HD23	1.98	0.45
1:D:385[A]:ARG:NH2	1:D:389:ARG:HH12	2.14	0.45
1:B:403:GLU:O	1:B:406[A]:GLN:CD	2.59	0.45
1:D:113:CYS:HB3	1:D:116:ILE:HD12	1.98	0.45
2:H:142:GLU:OE1	2:H:171:ARG:HD2	2.17	0.45
1:C:386:GLU:OE1	1:C:389:ARG:HD2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:133:ASP:OD1	2:E:133:ASP:N	2.44	0.44
1:A:164:LEU:HD12	1:A:165:PRO:HD2	1.97	0.44
2:H:36:LYS:N	2:H:36:LYS:CD	2.76	0.44
2:F:84:VAL:O	2:F:100:LYS:NZ	2.40	0.44
1:A:345:PRO:O	14:A:606:HOH:O	2.21	0.44
1:D:100:THR:HG21	1:D:122:PHE:CD1	2.53	0.44
2:F:138:ASP:O	2:F:141:ASN:N	2.49	0.44
1:A:157:ILE:HG13	1:A:158:PRO:HD2	1.98	0.44
1:B:29:ARG:NH1	14:B:629:HOH:O	2.50	0.44
2:G:112:ASN:HB2	2:G:116:ALA:HB3	2.00	0.44
2:G:126:ILE:O	2:G:130:ILE:HG12	2.18	0.44
1:C:106:THR:HG22	2:F:187[B]:LYS:HG3	1.99	0.44
2:H:101:ASP:OD1	2:H:101:ASP:C	2.60	0.44
2:H:133:ASP:OD1	2:H:133:ASP:N	2.38	0.44
2:F:142:GLU:OE1	2:F:171:ARG:HG2	2.17	0.44
2:H:127:LEU:O	2:H:131:GLY:N	2.50	0.44
1:C:52:ILE:HG12	4:C:502[A]:PG4:H22	2.00	0.44
2:F:69:GLU:OE1	2:F:96:LYS:HE2	2.18	0.44
2:G:164:LEU:HD13	2:H:32:TYR:HB3	2.00	0.44
2:G:208:PRO:O	2:G:212:LEU:HD23	2.18	0.44
2:H:86:TYR:CD2	2:H:195:PRO:HB3	2.52	0.44
2:G:140:ILE:C	2:G:142:GLU:H	2.26	0.43
2:F:36:LYS:HD2	2:F:36:LYS:HA	1.85	0.43
1:B:53:HIS:HB3	1:B:129:PRO:HG2	2.00	0.43
1:B:223:HIS:ND1	7:B:503:BU3:C3	2.81	0.43
1:C:174:LYS:HE3	14:C:1090:HOH:O	2.17	0.43
2:E:105:ARG:CZ	14:E:425:HOH:O	2.66	0.43
2:F:206:LYS:HG3	2:F:228:PRO:HG2	2.00	0.43
1:A:171:ARG:NH2	14:A:614:HOH:O	2.51	0.43
1:B:216:ARG:NH2	1:B:257:ASP:OD1	2.44	0.43
2:E:31:ARG:NH1	14:E:409:HOH:O	2.32	0.43
1:A:74:PHE:HE1	1:A:137:VAL:HG13	1.82	0.43
1:A:201:ARG:HD2	14:A:872:HOH:O	2.17	0.43
2:H:3:ALA:HB3	2:H:51:LEU:HD13	1.99	0.43
2:H:13:LYS:NZ	8:H:301:ADP:O1B	2.49	0.43
1:B:24:VAL:HB	1:B:25:PRO:HD3	2.00	0.43
1:C:421:LYS:NZ	14:C:625:HOH:O	2.40	0.43
2:E:18:LYS:HD2	14:E:414:HOH:O	2.19	0.43
2:F:148:ILE:HD12	2:F:177:LEU:CD2	2.49	0.43
1:B:22:ALA:HB3	1:B:23:PRO:HD3	2.01	0.43
1:B:95:GLY:O	1:B:285:PRO:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:138:ASP:C	2:H:171:ARG:HH12	2.20	0.43
1:D:215:LYS:HB3	1:D:215:LYS:HE3	1.80	0.43
2:E:206:LYS:O	2:E:208:PRO:HD3	2.19	0.43
1:A:140:GLU:HG2	1:A:142:THR:HG23	2.01	0.43
2:E:71:THR:HG23	2:E:105:ARG:NH1	2.32	0.42
2:F:84:VAL:CG1	2:F:199:ILE:HG13	2.48	0.42
1:B:290:LYS:HD2	1:B:406[A]:GLN:OE1	2.19	0.42
1:A:404:GLU:OE1	14:A:607:HOH:O	2.22	0.42
2:H:164:LEU:CD1	2:H:173:VAL:HG21	2.49	0.42
1:C:66:GLU:HG2	14:C:906:HOH:O	2.19	0.42
1:C:93:LEU:HB3	4:C:502[B]:PG4:H71	2.01	0.42
1:C:68:LYS:HE2	1:C:87:GLY:C	2.44	0.42
2:G:1:MET:HE1	2:G:19:GLU:CB	2.47	0.42
2:E:142:GLU:OE2	2:E:143:GLU:N	2.52	0.42
1:B:169:GLU:O	1:B:173[A]:ARG:HG3	2.19	0.42
2:H:160:VAL:O	2:H:164:LEU:HD13	2.20	0.42
1:D:64:LYS:HE3	1:D:135:ASP:OD2	2.20	0.42
2:E:157:GLU:OE2	2:E:161:ILE:HD11	2.20	0.42
1:A:75:CYS:SG	1:A:77:TYR:HB2	2.59	0.42
1:A:390:ARG:NE	14:A:639:HOH:O	2.52	0.42
1:B:164:LEU:HD12	1:B:165:PRO:HD2	2.02	0.42
2:H:100:LYS:HD2	2:H:104:GLY:HA2	2.01	0.42
1:D:29[A]:ARG:HA	1:D:33:LEU:HB2	2.02	0.42
2:E:48:ILE:O	14:E:404:HOH:O	2.22	0.42
2:H:97:ILE:HD13	2:H:189:ALA:HA	2.01	0.42
2:F:146:VAL:HG23	2:F:146:VAL:O	2.20	0.42
1:B:1:MET:HE3	1:B:1:MET:HA	2.01	0.42
1:B:380:SER:CB	1:B:396:LYS:HZ3	2.32	0.42
2:G:105:ARG:NE	14:G:421:HOH:O	2.53	0.42
2:E:43:LYS:HD3	2:E:43:LYS:HA	1.84	0.41
2:F:19:GLU:O	2:F:20:ASP:C	2.63	0.41
1:A:115:LEU:HD23	1:A:115:LEU:C	2.44	0.41
1:B:180:ILE:HG13	1:B:332:ALA:CB	2.50	0.41
2:H:164:LEU:N	2:H:164:LEU:CD1	2.82	0.41
2:H:126:ILE:CG2	2:H:127:LEU:N	2.83	0.41
1:D:16:LYS:HB3	1:D:127:ILE:CD1	2.50	0.41
2:E:105:ARG:CZ	14:E:403:HOH:O	2.62	0.41
2:F:130:ILE:HG21	2:F:170:LYS:HB3	2.02	0.41
2:H:65:LYS:NZ	2:H:243:THR:HG21	2.35	0.41
2:H:139:LEU:HD23	2:H:139:LEU:HA	1.83	0.41
2:H:184:ARG:NH1	14:H:421:HOH:O	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:17:ILE:HD11	2:F:234:LEU:HD12	2.03	0.41
2:H:14:LEU:O	2:H:25:PHE:HA	2.19	0.41
2:E:89:ASP:C	2:E:89:ASP:OD1	2.63	0.41
2:F:138:ASP:C	2:F:171:ARG:HH11	2.26	0.41
1:C:114:PRO:HB2	1:C:401:PHE:CE2	2.56	0.41
1:D:75:CYS:SG	1:D:77:TYR:HB2	2.60	0.41
2:E:148:ILE:H	2:E:180[B]:THR:CG2	2.33	0.41
2:F:148:ILE:HD12	2:F:177:LEU:HG	2.02	0.41
1:A:125:GLU:HB2	1:A:127:ILE:CD1	2.50	0.41
2:H:211:ARG:HD2	2:H:224:VAL:HG21	2.02	0.41
1:B:4:ARG:NH2	1:B:18:ASP:CG	2.77	0.41
1:B:385[B]:ARG:NH1	1:B:389:ARG:HH12	2.18	0.41
2:H:164:LEU:HD11	2:H:173:VAL:HG21	2.03	0.41
2:G:43:LYS:O	2:G:44:ASN:HB2	2.20	0.41
2:F:100:LYS:HA	2:F:105:ARG:O	2.21	0.41
1:A:382:LEU:HG	14:A:811:HOH:O	2.20	0.41
1:B:290:LYS:O	1:B:294:LEU:HG	2.21	0.41
1:C:29:ARG:HD2	1:C:34:ASN:CG	2.47	0.40
1:D:54:GLY:O	1:D:55:ILE:C	2.64	0.40
1:A:29:ARG:HD2	1:A:34:ASN:OD1	2.21	0.40
1:C:355:SER:OG	1:C:359:LYS:HE2	2.22	0.40
2:F:140:ILE:N	2:F:140:ILE:CD1	2.84	0.40
2:H:136:ASP:OD1	2:H:209:ALA:CA	2.68	0.40
1:D:98:GLN:NE2	14:E:413:HOH:O	2.54	0.40
2:F:39:GLU:HB3	14:F:401:HOH:O	2.20	0.40
1:B:282:THR:OG1	1:B:369:SER:HA	2.21	0.40
2:H:89:ASP:OD1	2:H:89:ASP:C	2.64	0.40
2:H:106:VAL:HG23	2:H:106:VAL:O	2.21	0.40
1:C:29:ARG:HD2	1:C:34:ASN:OD1	2.22	0.40
2:F:21:GLY:HA2	2:F:241:ARG:HH22	1.84	0.40
2:E:137:GLU:H	2:E:137:GLU:HG2	1.73	0.40
2:H:49:ARG:O	2:H:50:ASN:OD1	2.40	0.40

All (7) 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)
14:C:655:HOH:O	14:E:438:HOH:O[1_565]	1.95	0.25
14:A:975:HOH:O	14:G:428:HOH:O[1_545]	2.02	0.18
14:C:1170:HOH:O	14:A:1105:HOH:O[1_565]	2.06	0.14

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:A:605:HOH:O	14:G:413:HOH:O[1_545]	2.07	0.13
1:C:66:GLU:O	14:E:401:HOH:O[1_565]	2.15	0.05
14:C:955:HOH:O	14:G:474:HOH:O[1_655]	2.15	0.05
14:A:1120:HOH:O	14:G:413:HOH:O[1_545]	2.15	0.05

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	425/422 (101%)	415 (98%)	10 (2%)	0	100	100
1	B	430/422 (102%)	420 (98%)	10 (2%)	0	100	100
1	C	421/422 (100%)	410 (97%)	11 (3%)	0	100	100
1	D	426/422 (101%)	416 (98%)	10 (2%)	0	100	100
2	E	245/243 (101%)	240 (98%)	3 (1%)	2 (1%)	16	8
2	F	243/243 (100%)	238 (98%)	4 (2%)	1 (0%)	30	21
2	G	241/243 (99%)	233 (97%)	8 (3%)	0	100	100
2	H	242/243 (100%)	240 (99%)	2 (1%)	0	100	100
All	All	2673/2660 (100%)	2612 (98%)	58 (2%)	3 (0%)	48	41

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	136	ASP
2	E	131	GLY
2	F	131	GLY

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	369/363 (102%)	367 (100%)	2 (0%)	81	82
1	B	373/363 (103%)	371 (100%)	2 (0%)	81	82
1	C	365/363 (101%)	365 (100%)	0	100	100
1	D	370/363 (102%)	368 (100%)	2 (0%)	81	82
2	E	200/196 (102%)	198 (99%)	2 (1%)	68	67
2	F	198/196 (101%)	196 (99%)	2 (1%)	68	67
2	G	196/196 (100%)	196 (100%)	0	100	100
2	H	197/196 (100%)	197 (100%)	0	100	100
All	All	2268/2236 (101%)	2258 (100%)	10 (0%)	87	85

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	385[A]	ARG
1	D	385[B]	ARG
2	E	211[A]	ARG
2	E	211[B]	ARG
2	F	194[A]	LYS
2	F	194[B]	LYS
1	A	19[A]	GLU
1	A	19[B]	GLU
1	B	123[A]	LYS
1	B	123[C]	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	361	GLN
1	D	62	ASN
2	E	103	GLN
2	E	141	ASN
2	F	73	GLN
2	F	144	ASN
2	F	215	GLN
1	A	65	GLN

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Mol	Chain	Res	Type
1	A	132	GLN
1	A	371	GLN
1	B	11	ASN
1	B	37	ASN
1	B	62	ASN
1	B	361	GLN
2	G	44	ASN
2	H	11	ASN
2	H	141	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](#)

Of 36 ligands modelled in this entry, 4 are monoatomic and 2 are modelled with single atom - leaving 30 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	$\# Z > 2$
6	TAM	B	504	-	10,10,10	0.54	0	12,12,12	1.71	4 (33%)
10	ALF	G	303	-	4,4,4	1.41	0	-		
12	SO4	A	501	-	4,4,4	0.30	0	6,6,6	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BJ8	B	501	1	6,26,26	4.27	4 (66%)	-		
3	BJ8	C	501	1	6,26,26	3.86	4 (66%)	-		
6	TAM	E	304	-	10,10,10	0.45	0	12,12,12	1.22	2 (16%)
7	BU3	G	307	-	4,5,5	0.51	0	6,6,6	0.63	0
10	ALF	E	303	-	4,4,4	1.36	0	-		
12	SO4	A	503	-	4,4,4	0.34	0	6,6,6	0.15	0
6	TAM	G	306	-	10,10,10	0.76	0	12,12,12	2.57	3 (25%)
8	ADP	H	301	9	28,29,29	1.61	6 (21%)	43,45,45	1.98	9 (20%)
12	SO4	A	505	-	4,4,4	0.27	0	6,6,6	0.18	0
4	PG4	C	502[B]	-	12,12,12	0.55	0	11,11,11	0.49	0
6	TAM	D	503	-	10,10,10	0.68	0	12,12,12	1.32	1 (8%)
12	SO4	G	304	-	4,4,4	0.38	0	6,6,6	0.24	0
12	SO4	F	305	-	4,4,4	0.25	0	6,6,6	0.29	0
4	PG4	C	502[A]	-	12,12,12	0.65	0	11,11,11	0.70	0
11	SF4	F	301	2	0,12,12	-	-	-		
7	BU3	B	503	-	4,5,5	0.66	0	6,6,6	0.77	0
10	ALF	F	304	-	4,4,4	1.49	0	-		
8	ADP	E	301	9	28,29,29	1.75	5 (17%)	43,45,45	1.97	13 (30%)
7	BU3	D	504	-	4,5,5	0.69	0	6,6,6	0.41	0
3	BJ8	D	501	1	6,26,26	4.13	4 (66%)	-		
3	BJ8	A	504	1	6,26,26	3.53	3 (50%)	-		
7	BU3	B	502	-	4,5,5	0.82	0	6,6,6	0.36	0
8	ADP	F	302	9	28,29,29	1.41	4 (14%)	43,45,45	1.90	14 (32%)
8	ADP	G	301	9	28,29,29	1.68	5 (17%)	43,45,45	1.93	11 (25%)
11	SF4	G	308	2	0,12,12	-	-	-		
5	GOL	D	502	-	5,5,5	0.99	0	5,5,5	0.89	0
10	ALF	H	303	-	4,4,4	1.46	0	-		

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
6	TAM	B	504	-	-	8/12/12/12	-
3	BJ8	B	501	1	-	-	0/12/10/10
3	BJ8	C	501	1	-	-	0/12/10/10
6	TAM	E	304	-	-	6/12/12/12	-
7	BU3	G	307	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	TAM	G	306	-	-	3/12/12/12	-
8	ADP	H	301	9	-	2/16/32/32	0/3/3/3
4	PG4	C	502[B]	-	-	4/10/10/10	-
6	TAM	D	503	-	-	9/12/12/12	-
4	PG4	C	502[A]	-	-	4/10/10/10	-
11	SF4	F	301	2	-	-	0/6/5/5
7	BU3	B	503	-	-	1/4/4/4	-
8	ADP	E	301	9	-	1/16/32/32	0/3/3/3
7	BU3	D	504	-	-	0/4/4/4	-
3	BJ8	D	501	1	-	-	0/12/10/10
3	BJ8	A	504	1	-	-	0/12/10/10
7	BU3	B	502	-	-	0/4/4/4	-
8	ADP	F	302	9	-	3/16/32/32	0/3/3/3
8	ADP	G	301	9	-	1/16/32/32	0/3/3/3
11	SF4	G	308	2	-	-	0/6/5/5
5	GOL	D	502	-	-	4/4/4/4	-

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	BJ8	S5-FE8	8.25	2.41	2.28
3	C	501	BJ8	S5-FE8	7.10	2.39	2.28
3	D	501	BJ8	S6-FE8	6.73	2.38	2.28
3	D	501	BJ8	S5-FE8	5.49	2.36	2.28
8	H	301	ADP	C5-C4	5.48	1.48	1.39
8	E	301	ADP	C5-C4	5.47	1.48	1.39
8	G	301	ADP	C5-C4	5.40	1.48	1.39
3	A	504	BJ8	S7-FE8	4.88	2.35	2.28
3	A	504	BJ8	S5-FE8	4.76	2.35	2.28
8	E	301	ADP	PA-O3A	4.61	1.64	1.59
3	A	504	BJ8	S2-FE1	4.55	2.35	2.28
8	F	302	ADP	C5-C4	4.39	1.46	1.39
3	C	501	BJ8	S6-FE8	4.33	2.34	2.28
3	D	501	BJ8	S2-FE1	4.10	2.34	2.28
3	B	501	BJ8	S6-FE8	4.06	2.34	2.28
3	B	501	BJ8	S2-FE1	4.05	2.34	2.28
3	C	501	BJ8	S3-FE1	3.91	2.34	2.28
8	G	301	ADP	PA-O3A	3.31	1.63	1.59
8	H	301	ADP	C4-N9	-3.10	1.31	1.37
8	F	302	ADP	C5-C6	2.89	1.49	1.41
8	H	301	ADP	C8-N7	2.84	1.37	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	501	BJ8	S7-FE8	-2.76	2.23	2.28
8	G	301	ADP	C5-C6	2.75	1.48	1.41
8	E	301	ADP	C5-N7	-2.71	1.34	1.39
3	B	501	BJ8	S4-FE1	-2.64	2.23	2.28
8	E	301	ADP	C5-C6	2.57	1.48	1.41
8	G	301	ADP	C4-N9	-2.53	1.32	1.37
8	F	302	ADP	C4-N9	-2.50	1.32	1.37
8	H	301	ADP	PA-O3A	2.44	1.62	1.59
8	H	301	ADP	C5-C6	2.31	1.47	1.41
8	H	301	ADP	C2-N3	2.22	1.37	1.33
8	G	301	ADP	C5-N7	-2.18	1.35	1.39
8	E	301	ADP	C4-N9	-2.09	1.33	1.37
8	F	302	ADP	C5-N7	-2.09	1.35	1.39
3	C	501	BJ8	S4-FE1	-2.01	2.24	2.28

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	G	306	TAM	C6-C3-C	-5.70	109.25	115.97
8	E	301	ADP	C5-C4-N3	-5.37	119.32	126.72
6	G	306	TAM	C4-C1-C	-5.19	109.84	115.97
8	G	301	ADP	C5-C4-N3	-5.04	119.78	126.72
8	F	302	ADP	C5-C4-N3	-5.03	119.79	126.72
8	H	301	ADP	C5-C4-N3	-4.88	120.00	126.72
8	G	301	ADP	N3-C4-N9	4.71	135.18	127.17
8	E	301	ADP	N3-C4-N9	4.61	135.00	127.17
8	H	301	ADP	N3-C2-N1	-4.41	121.91	128.58
8	E	301	ADP	N3-C2-N1	-4.32	122.05	128.58
8	H	301	ADP	N3-C4-N9	4.26	134.41	127.17
8	F	302	ADP	N3-C4-N9	4.13	134.20	127.17
8	G	301	ADP	N3-C2-N1	-3.99	122.54	128.58
6	G	306	TAM	C5-C2-C	-3.96	111.30	115.97
8	E	301	ADP	C2-N3-C4	3.86	121.26	111.83
8	F	302	ADP	C3'-C2'-C1'	3.83	108.71	101.46
8	H	301	ADP	C4-N9-C8	3.77	109.70	105.74
8	F	302	ADP	N3-C2-N1	-3.69	123.00	128.58
8	H	301	ADP	C3'-C2'-C1'	3.60	108.27	101.46
8	H	301	ADP	C2-N3-C4	3.57	120.56	111.83
8	G	301	ADP	C4-N9-C8	3.51	109.42	105.74
8	G	301	ADP	C2-N3-C4	3.44	120.23	111.83
8	F	302	ADP	C2-N3-C4	3.44	120.23	111.83
6	B	504	TAM	C2-C-N	3.42	116.81	108.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	F	302	ADP	C4-C5-N7	-3.31	106.80	110.58
8	E	301	ADP	C3'-C2'-C1'	3.19	107.50	101.46
6	D	503	TAM	C3-C-C1	-3.16	104.94	110.50
6	B	504	TAM	C2-C-C1	-3.12	105.01	110.50
8	F	302	ADP	C4-N9-C8	3.07	108.96	105.74
8	G	301	ADP	C4-C5-N7	-2.94	107.22	110.58
8	E	301	ADP	C4-C5-N7	-2.87	107.30	110.58
8	H	301	ADP	C2-N1-C6	2.81	123.34	118.73
8	G	301	ADP	C3'-C2'-C1'	2.78	106.72	101.46
8	G	301	ADP	O3A-PB-O1B	-2.77	96.47	111.04
8	E	301	ADP	C2-N1-C6	2.64	123.07	118.73
8	E	301	ADP	C5-N7-C8	2.63	107.58	103.45
8	H	301	ADP	O4'-C1'-N9	2.61	113.11	108.09
8	F	302	ADP	C2-N1-C6	2.55	122.93	118.73
6	E	304	TAM	C6-C3-C	-2.53	112.98	115.97
8	G	301	ADP	O2B-PB-O3A	2.50	113.03	104.64
6	E	304	TAM	C4-C1-C	-2.49	113.03	115.97
6	B	504	TAM	C6-C3-C	-2.49	113.03	115.97
8	H	301	ADP	C4-C5-N7	-2.49	107.74	110.58
8	E	301	ADP	C4-N9-C8	2.49	108.35	105.74
8	G	301	ADP	C2-N1-C6	2.44	122.75	118.73
8	E	301	ADP	C6-C5-N7	2.33	136.59	132.09
8	E	301	ADP	O4'-C1'-N9	2.33	112.56	108.09
8	F	302	ADP	C5-N7-C8	2.30	107.07	103.45
8	G	301	ADP	C5-N7-C8	2.21	106.92	103.45
8	F	302	ADP	O2A-PA-O1A	2.17	122.55	112.44
8	F	302	ADP	C6-C5-N7	2.16	136.25	132.09
8	F	302	ADP	O4'-C1'-N9	2.14	112.20	108.09
8	E	301	ADP	O3B-PB-O2B	2.09	115.65	107.80
8	F	302	ADP	O3A-PB-O1B	-2.09	100.06	111.04
8	F	302	ADP	O3B-PB-O2B	2.07	115.55	107.80
8	E	301	ADP	O2A-PA-O1A	2.06	122.03	112.44
6	B	504	TAM	C3-C-C2	-2.05	106.89	110.50

There are no chirality outliers.

All (50) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	502	GOL	O1-C1-C2-C3
5	D	502	GOL	O2-C2-C3-O3
6	D	503	TAM	C2-C-C1-C4
6	D	503	TAM	C3-C-C1-C4

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Mol	Chain	Res	Type	Atoms
6	D	503	TAM	N-C-C1-C4
6	D	503	TAM	C3-C-C2-C5
6	D	503	TAM	N-C-C2-C5
6	D	503	TAM	C2-C-C3-C6
6	D	503	TAM	N-C-C3-C6
6	E	304	TAM	N-C-C1-C4
6	B	504	TAM	N-C-C1-C4
6	B	504	TAM	C1-C-C2-C5
6	B	504	TAM	C3-C-C2-C5
6	B	504	TAM	N-C-C2-C5
6	B	504	TAM	C-C3-C6-O6
6	G	306	TAM	C1-C-C3-C6
6	G	306	TAM	C2-C-C3-C6
6	G	306	TAM	N-C-C3-C6
7	B	503	BU3	C1-C2-C3-C4
7	G	307	BU3	O5-C2-C3-O6
7	G	307	BU3	C1-C2-C3-O6
7	G	307	BU3	O5-C2-C3-C4
7	G	307	BU3	C1-C2-C3-C4
5	D	502	GOL	C1-C2-C3-O3
5	D	502	GOL	O1-C1-C2-O2
6	E	304	TAM	C-C2-C5-O5
6	E	304	TAM	C-C3-C6-O6
6	B	504	TAM	C-C1-C4-O4
4	C	502[B]	PG4	O1-C1-C2-O2
6	D	503	TAM	C1-C-C2-C5
6	D	503	TAM	C1-C-C3-C6
6	E	304	TAM	C3-C-C1-C4
8	F	302	ADP	PA-O3A-PB-O1B
4	C	502[B]	PG4	O2-C3-C4-O3
8	H	301	ADP	PA-O3A-PB-O3B
4	C	502[B]	PG4	C6-C5-O3-C4
6	B	504	TAM	C2-C-C1-C4
6	B	504	TAM	C3-C-C1-C4
4	C	502[A]	PG4	C6-C5-O3-C4
4	C	502[B]	PG4	C4-C3-O2-C2
6	E	304	TAM	C-C1-C4-O4
4	C	502[A]	PG4	C3-C4-O3-C5
4	C	502[A]	PG4	C4-C3-O2-C2
6	E	304	TAM	C2-C-C1-C4
8	G	301	ADP	O4'-C1'-N9-C8
8	F	302	ADP	C2'-C1'-N9-C4

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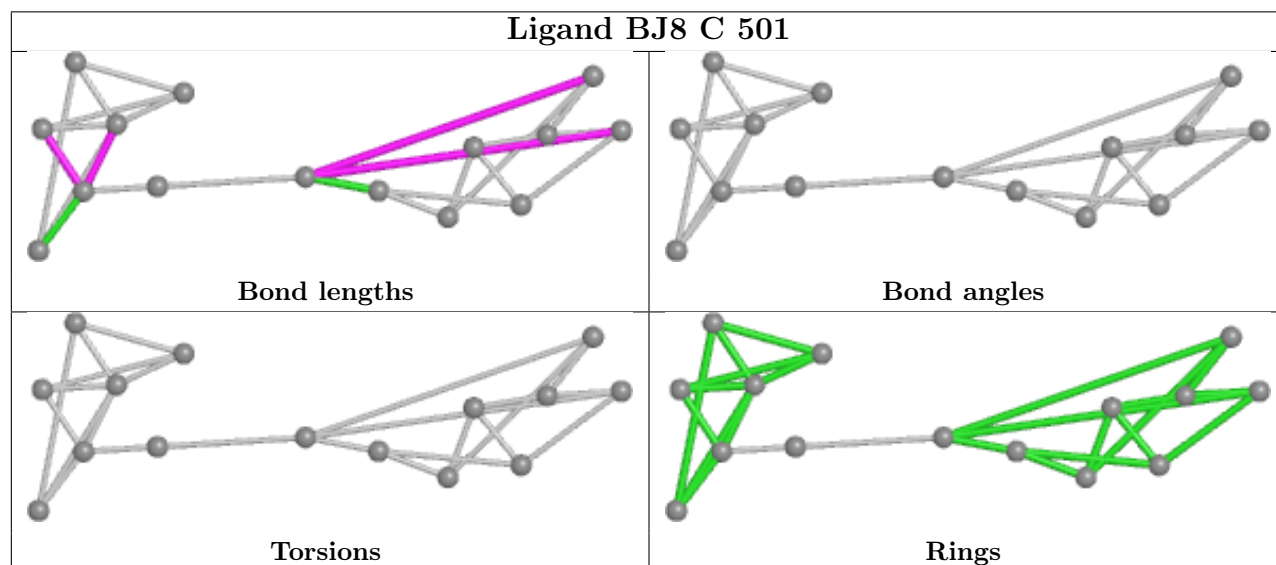
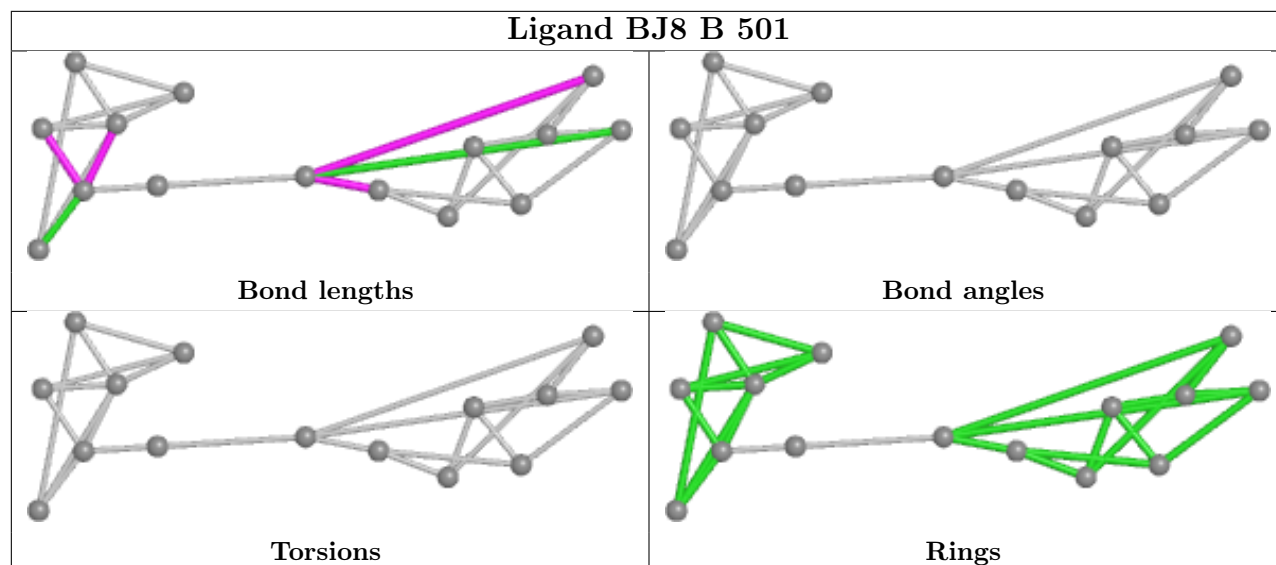
Mol	Chain	Res	Type	Atoms
8	E	301	ADP	O4'-C1'-N9-C8
8	F	302	ADP	O4'-C1'-N9-C8
8	H	301	ADP	O4'-C1'-N9-C8
4	C	502[A]	PG4	O4-C7-C8-O5

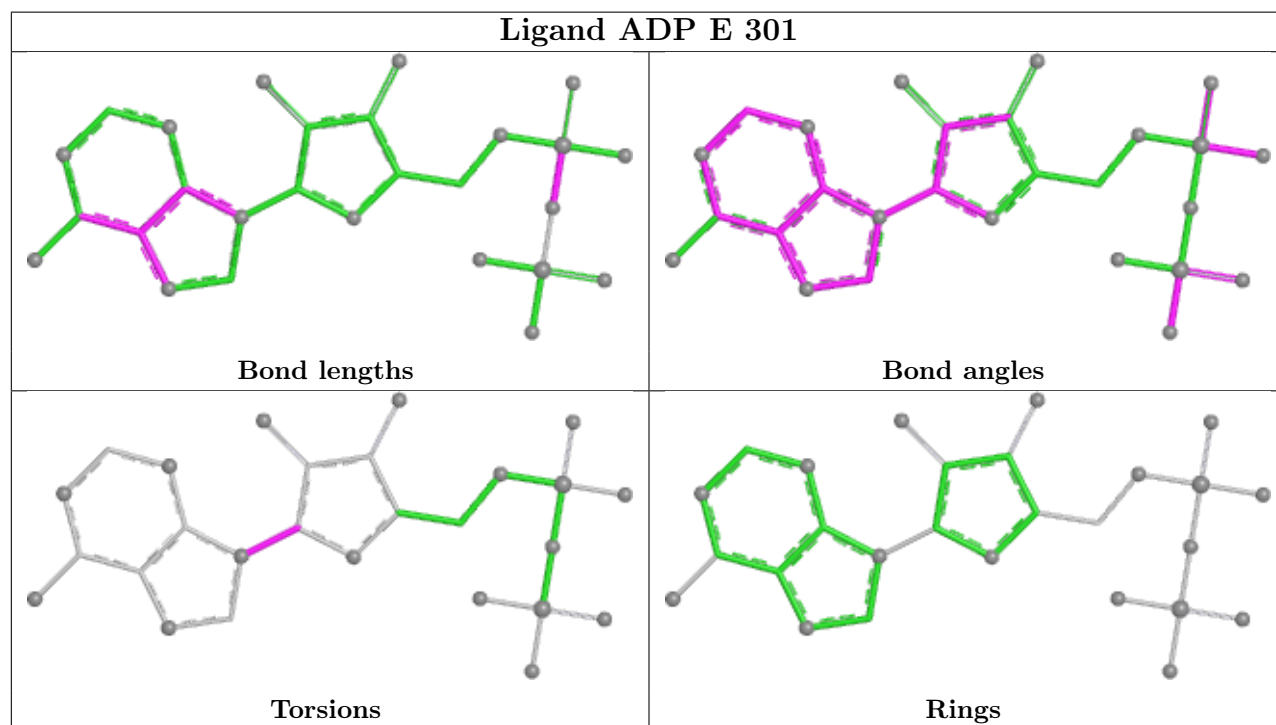
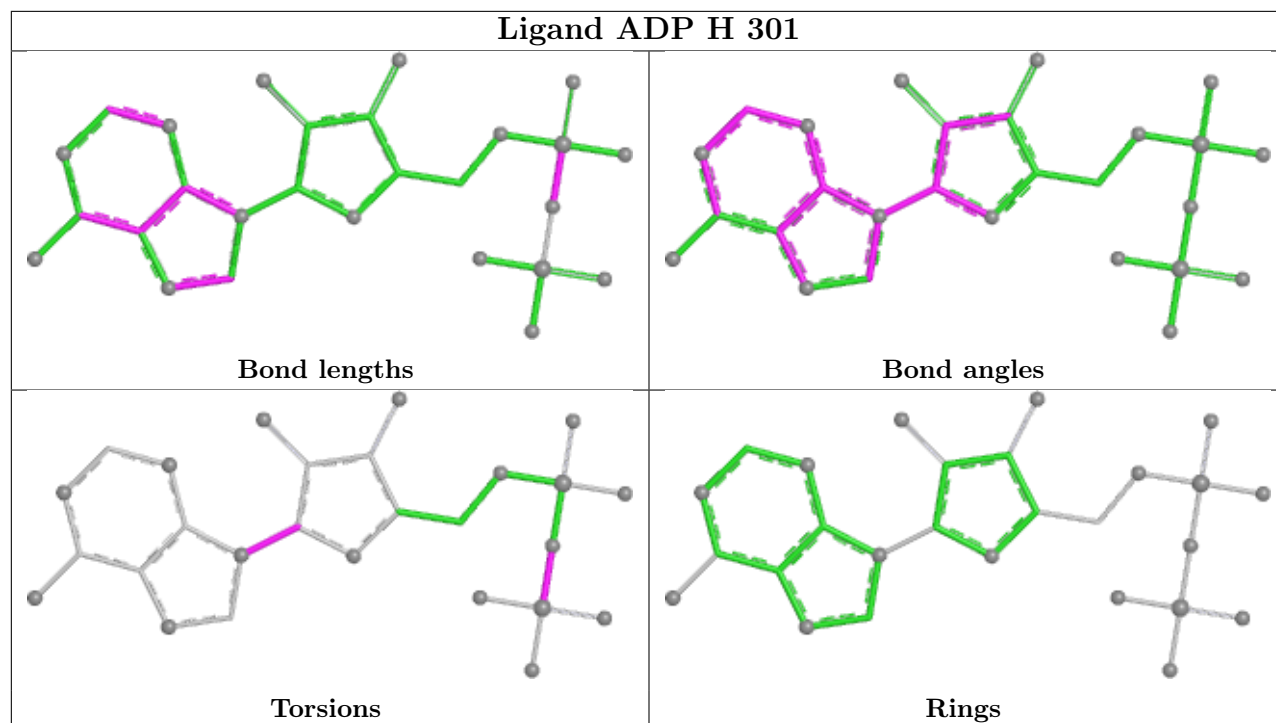
There are no ring outliers.

8 monomers are involved in 13 short contacts:

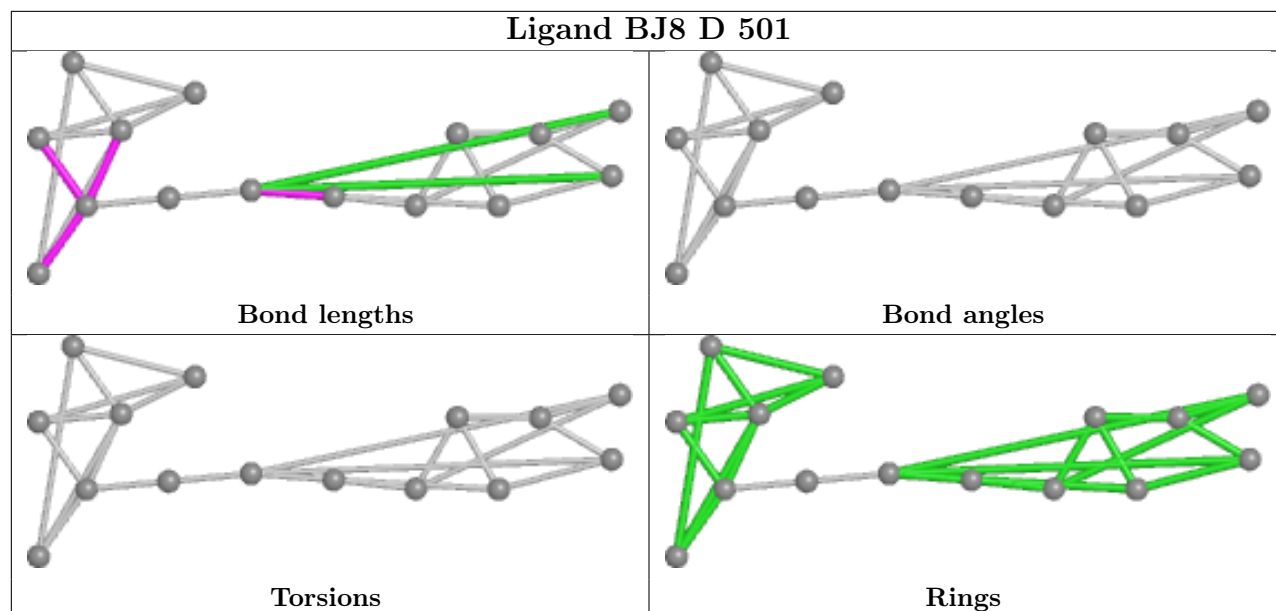
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	504	TAM	1	0
12	A	501	SO4	1	0
3	C	501	BJ8	1	0
8	H	301	ADP	1	0
4	C	502[B]	PG4	1	0
12	G	304	SO4	1	0
4	C	502[A]	PG4	4	0
7	B	503	BU3	3	0

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

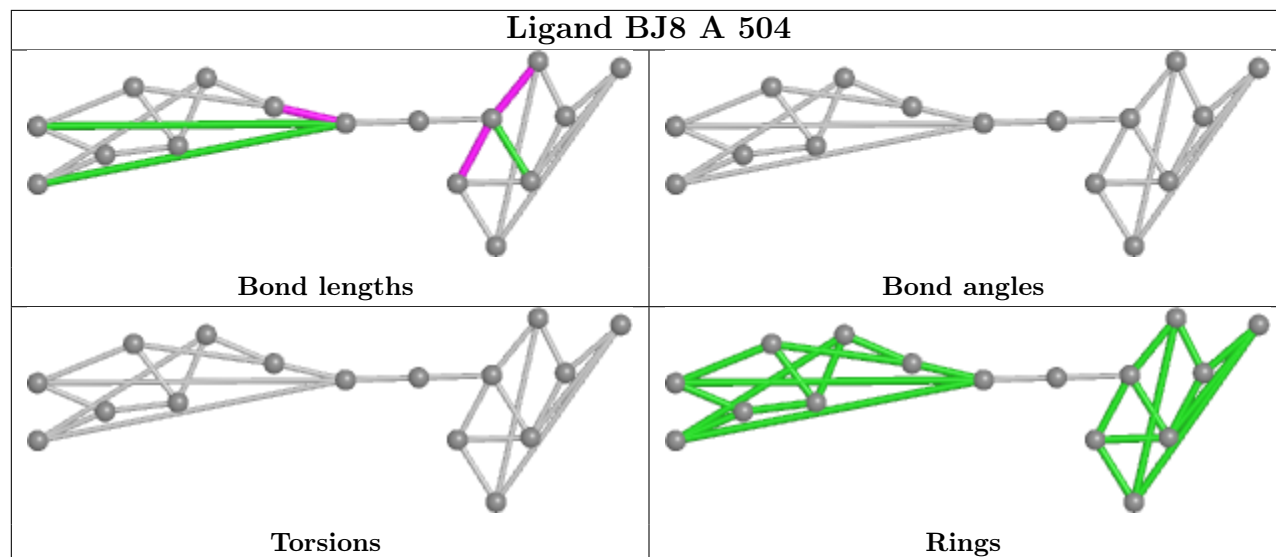


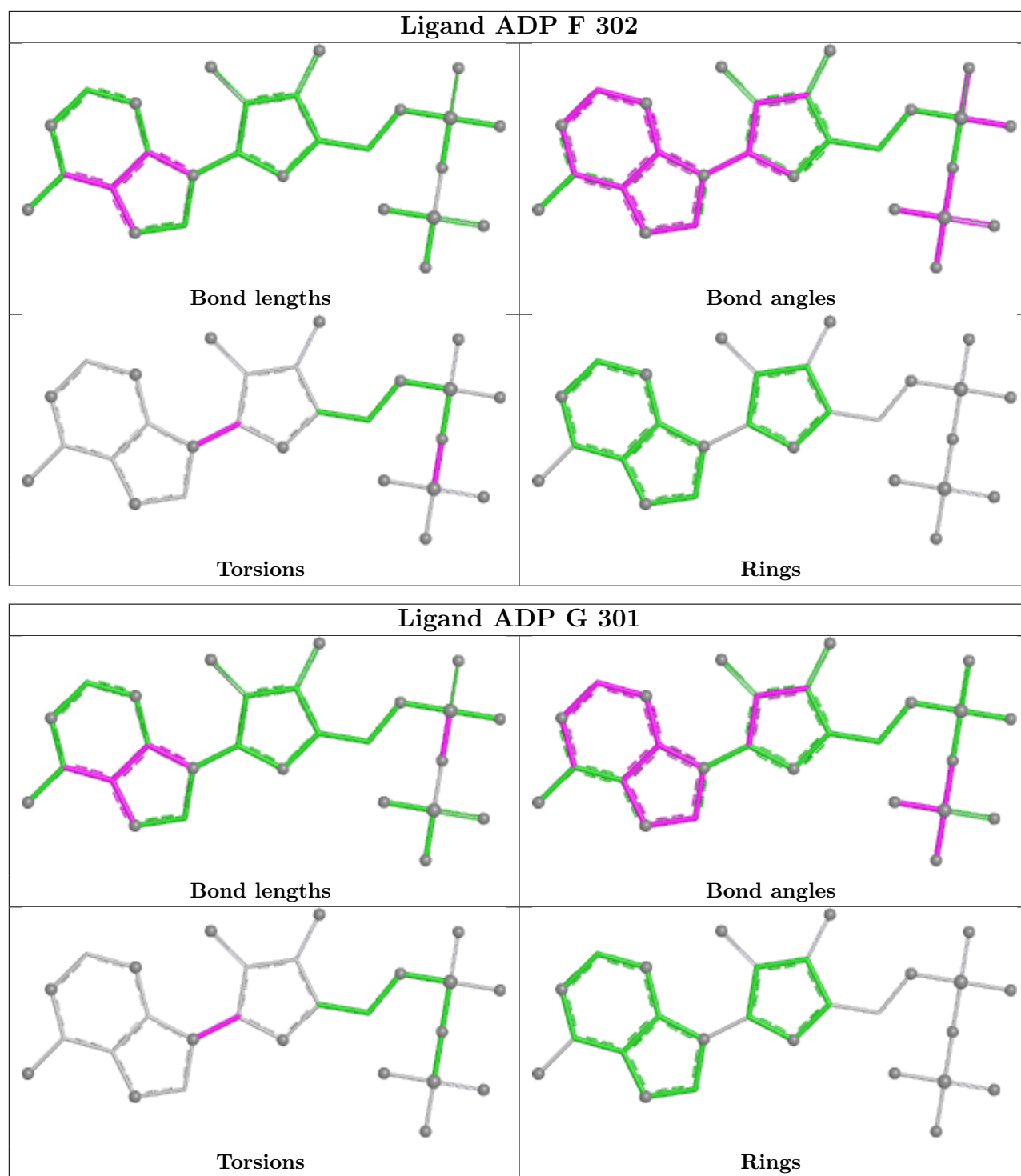


Ligand BJ8 D 501



Ligand BJ8 A 504





5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	421/422 (99%)	-0.50	2 (0%) 87 90	9, 19, 32, 87	6 (1%)
1	B	422/422 (100%)	-0.05	4 (0%) 81 85	13, 29, 44, 100	8 (1%)
1	C	420/422 (99%)	-0.53	0 100 100	10, 19, 32, 47	4 (0%)
1	D	421/422 (99%)	0.01	3 (0%) 84 88	9, 29, 44, 96	6 (1%)
2	E	243/243 (100%)	-0.04	16 (6%) 24 28	8, 23, 56, 121	13 (5%)
2	F	243/243 (100%)	0.74	22 (9%) 15 17	16, 36, 62, 84	15 (6%)
2	G	243/243 (100%)	-0.02	18 (7%) 20 23	12, 23, 60, 114	9 (3%)
2	H	243/243 (100%)	1.00	39 (16%) 5 5	15, 40, 75, 96	11 (4%)
All	All	2656/2660 (99%)	-0.01	104 (3%) 43 50	8, 26, 53, 121	72 (2%)

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	134	TYR	7.4
2	H	134	TYR	7.3
2	H	132	ASP	5.5
2	F	132	ASP	5.2
2	G	134	TYR	5.0
2	H	138	ASP	4.7
2	E	134	TYR	4.5
2	F	133	ASP	4.5
2	H	131	GLY	4.5
2	H	133	ASP	4.4
2	G	140	ILE	4.4
2	H	135	ARG	4.4
2	F	138	ASP	4.3
2	F	131	GLY	4.2
1	A	1	MET	4.1
2	F	136	ASP	4.0

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Mol	Chain	Res	Type	RSRZ
2	G	139	LEU	3.7
2	G	144	ASN	3.7
2	H	140	ILE	3.6
2	H	186	ALA	3.5
2	H	17	ILE	3.5
2	F	137	GLU	3.5
2	H	139	LEU	3.5
2	H	136	ASP	3.5
2	F	141	ASN	3.4
2	H	243	THR	3.4
2	E	144	ASN	3.4
2	F	140	ILE	3.4
2	G	136	ASP	3.3
1	B	1	MET	3.3
2	F	135	ARG	3.3
2	G	145	ALA	3.2
2	F	139	LEU	3.1
1	D	1	MET	3.1
2	E	131	GLY	3.1
2	H	74	ALA	3.1
2	H	81	PHE	3.0
2	E	133	ASP	3.0
2	E	139	LEU	3.0
2	E	19	GLU	3.0
2	H	102	GLY	3.0
2	E	145	ALA	3.0
2	H	141	ASN	2.9
2	F	48	ILE	2.9
2	H	42	LEU	2.9
2	E	138	ASP	2.9
2	H	50	ASN	2.9
2	E	136	ASP	2.9
1	B	0	GLY	2.9
2	F	243	THR	2.8
2	G	131	GLY	2.8
2	H	48	ILE	2.8
2	H	239	ILE	2.8
2	F	46	GLY	2.7
2	G	135	ARG	2.7
2	H	137	GLU	2.7
2	G	142	GLU	2.6
2	H	45	THR	2.6

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Mol	Chain	Res	Type	RSRZ
2	G	141	ASN	2.6
2	E	135	ARG	2.6
2	E	140	ILE	2.5
2	H	130	ILE	2.5
2	H	240	ALA	2.5
2	E	243	THR	2.4
2	E	141	ASN	2.4
2	G	32	TYR	2.4
2	H	49	ARG	2.4
2	F	59	VAL	2.4
2	H	82	PRO	2.4
1	D	421	LYS	2.4
2	G	243	THR	2.3
2	E	132	ASP	2.3
2	F	20	ASP	2.3
2	H	143	GLU	2.3
2	G	45	THR	2.3
2	H	2	PHE	2.2
1	B	55	ILE	2.2
2	F	44	ASN	2.2
2	F	82	PRO	2.2
2	H	46	GLY	2.2
2	F	81	PHE	2.2
1	B	421	LYS	2.2
2	H	1	MET	2.2
2	F	42	LEU	2.2
2	H	80	LEU	2.2
2	H	3	ALA	2.2
2	G	143	GLU	2.2
2	H	20	ASP	2.2
2	F	2	PHE	2.1
2	E	45	THR	2.1
2	E	48	ILE	2.1
1	A	19[A]	GLU	2.1
2	F	145	ALA	2.1
2	H	59	VAL	2.1
2	H	40	LEU	2.1
2	G	19	GLU	2.0
2	H	27	VAL	2.0
2	H	53	VAL	2.0
2	G	133	ASP	2.0
2	G	138	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	272	LYS	2.0
2	H	25	PHE	2.0
2	H	103	GLN	2.0
2	G	146	VAL	2.0

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

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

6.3 Carbohydrates ⓘ

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.

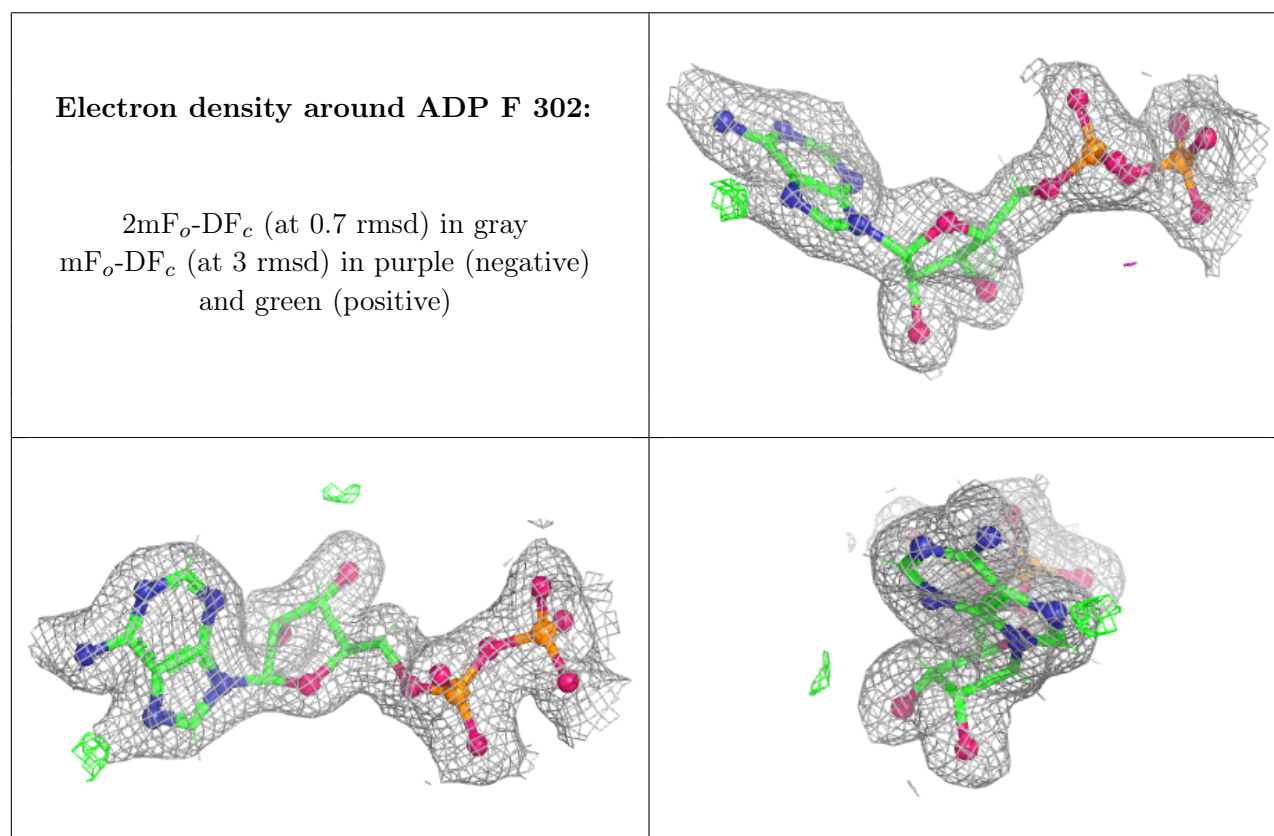
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	TAM	E	304	11/11	0.62	0.20	59,75,82,83	0
6	TAM	G	306	11/11	0.67	0.18	31,42,52,54	28
7	BU3	D	504	6/6	0.69	0.18	48,64,77,78	0
7	BU3	B	503	6/6	0.69	0.22	35,55,61,62	16
4	PG4	C	502[B]	13/13	0.73	0.16	26,32,39,42	31
4	PG4	C	502[A]	13/13	0.73	0.16	27,33,40,44	31
7	BU3	B	502	6/6	0.73	0.18	29,44,50,51	0
6	TAM	B	504	11/11	0.73	0.18	42,56,67,68	28
6	TAM	D	503	11/11	0.78	0.13	30,39,56,59	0
12	SO4	A	503	5/5	0.78	0.19	33,34,41,44	5
7	BU3	G	307	6/6	0.81	0.18	34,49,65,65	16
5	GOL	D	502	6/6	0.82	0.17	29,37,53,64	14
12	SO4	A	505	5/5	0.83	0.15	26,33,35,46	5
12	SO4	F	305	5/5	0.84	0.20	27,34,41,44	5
13	NH4	G	305	1/1	0.87	0.13	27,27,27,27	0
13	NH4	A	502	1/1	0.90	0.13	21,21,21,21	0
12	SO4	A	501	5/5	0.92	0.12	33,36,42,45	0
12	SO4	G	304	5/5	0.94	0.08	36,36,43,44	0
10	ALF	H	303	5/5	0.97	0.04	21,23,29,30	0

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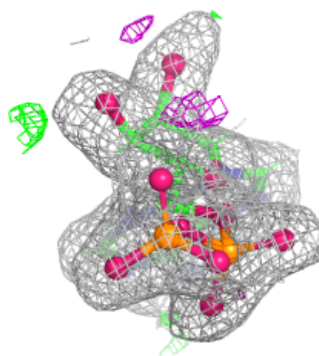
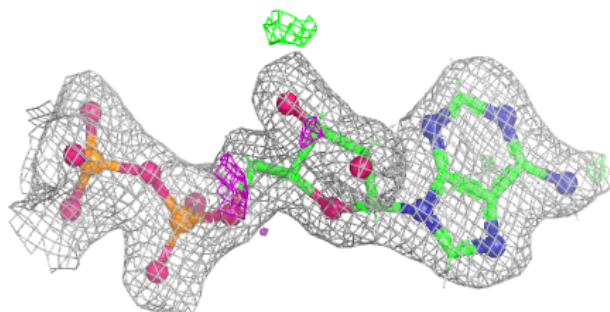
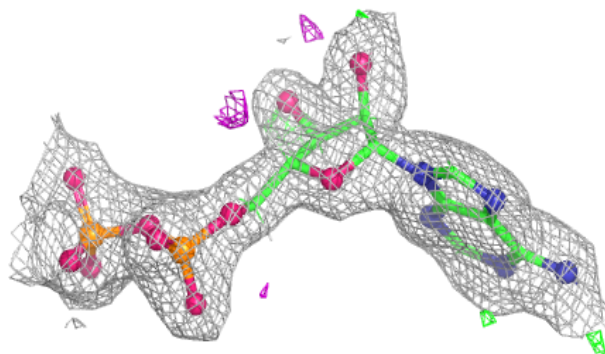
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	ADP	F	302	27/27	0.97	0.06	17,27,38,45	0
8	ADP	H	301	27/27	0.97	0.06	20,27,38,44	0
10	ALF	F	304	5/5	0.97	0.05	20,21,25,25	0
10	ALF	E	303	5/5	0.98	0.04	8,12,14,15	0
10	ALF	G	303	5/5	0.98	0.04	11,12,15,16	0
3	BJ8	A	504	17/17	0.99	0.03	12,15,18,19	0
8	ADP	E	301	27/27	0.99	0.05	10,19,38,46	0
3	BJ8	B	501	17/17	0.99	0.03	18,20,27,29	0
8	ADP	G	301	27/27	0.99	0.04	10,19,38,40	0
3	BJ8	C	501	17/17	0.99	0.03	11,14,18,18	0
9	MG	E	302	1/1	0.99	0.03	12,12,12,12	0
9	MG	F	303	1/1	0.99	0.03	21,21,21,21	0
9	MG	G	302	1/1	0.99	0.01	10,10,10,10	0
9	MG	H	302	1/1	0.99	0.03	25,25,25,25	0
3	BJ8	D	501	17/17	0.99	0.04	16,19,26,26	0
11	SF4	G	308	8/8	1.00	0.02	13,16,18,18	0
11	SF4	F	301	8/8	1.00	0.02	13,15,17,18	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.

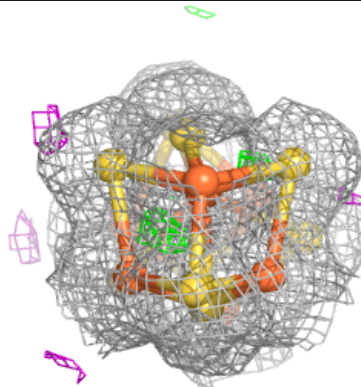
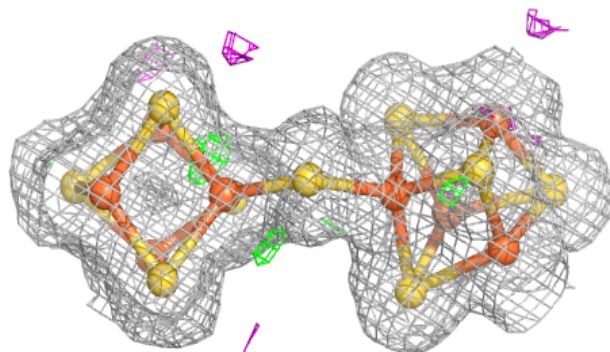
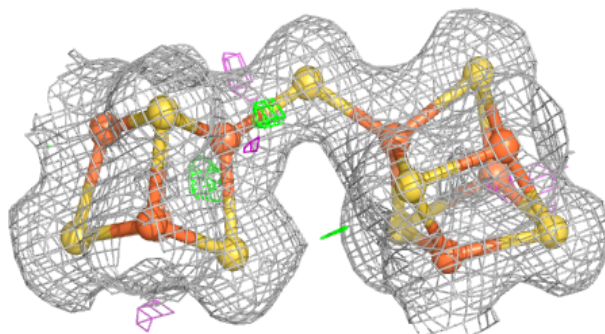


Electron density around ADP H 301:

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

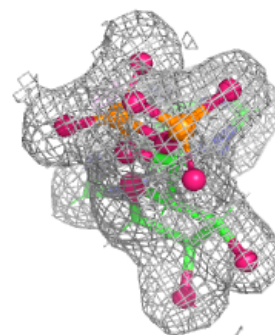
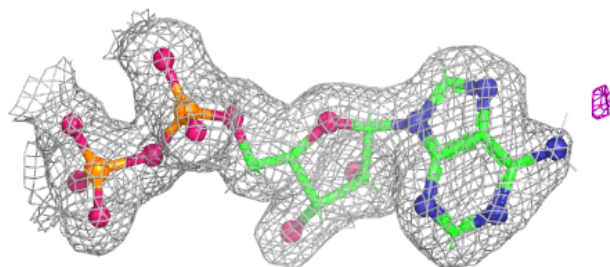
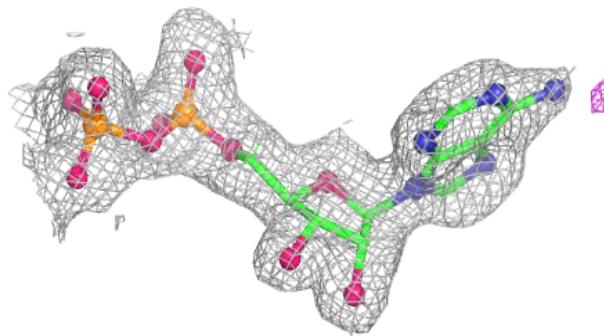
**Electron density around BJ8 A 504:**

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

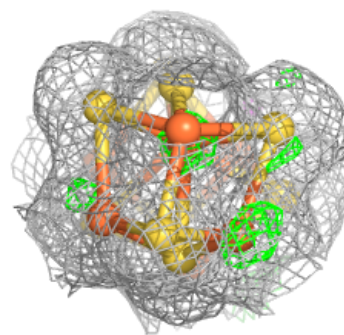
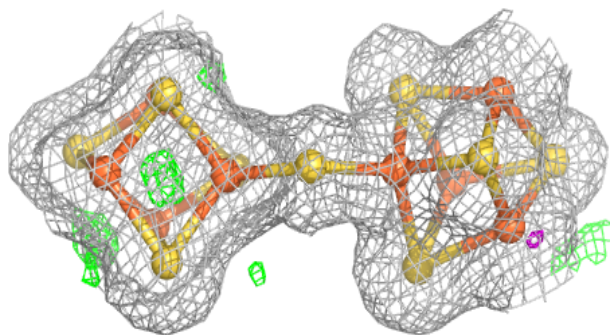
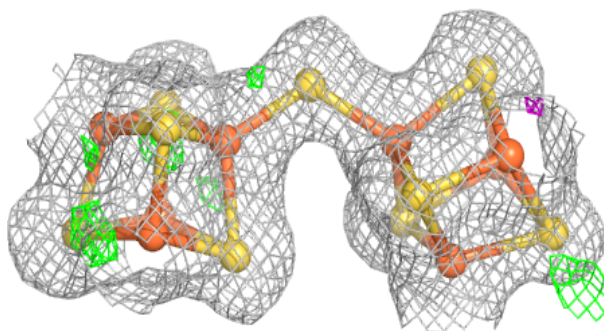


Electron density around ADP E 301:

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

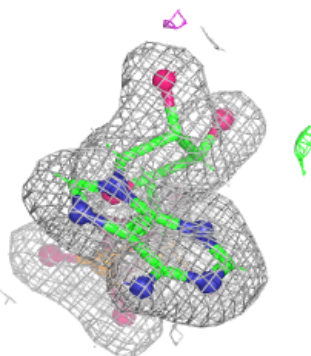
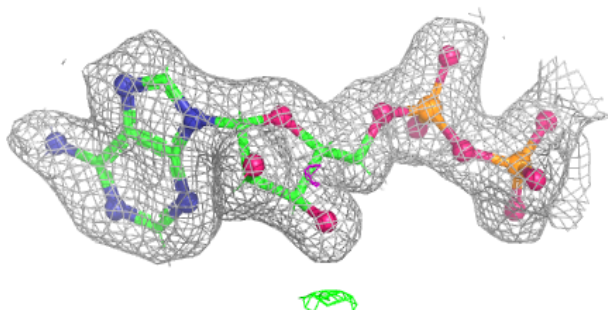
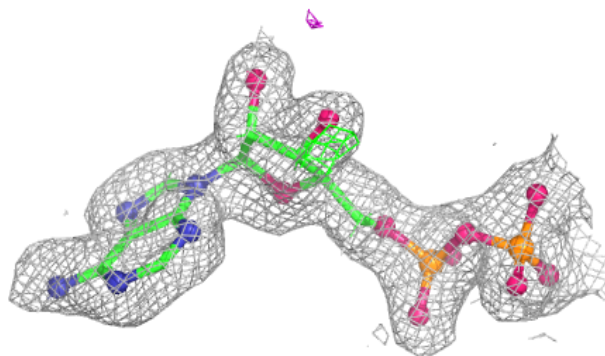
**Electron density around BJ8 B 501:**

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

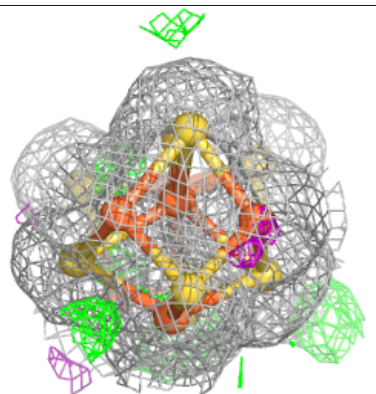
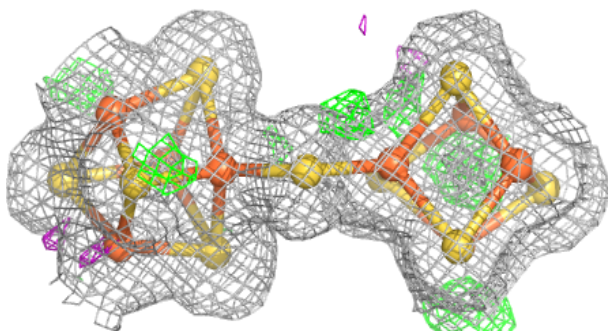
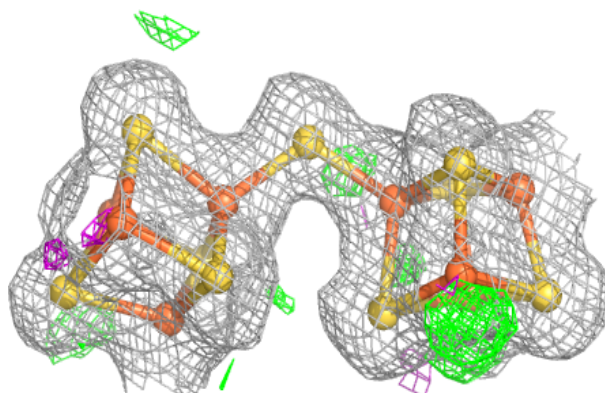


Electron density around ADP G 301:

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

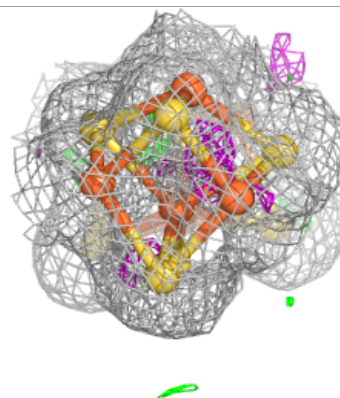
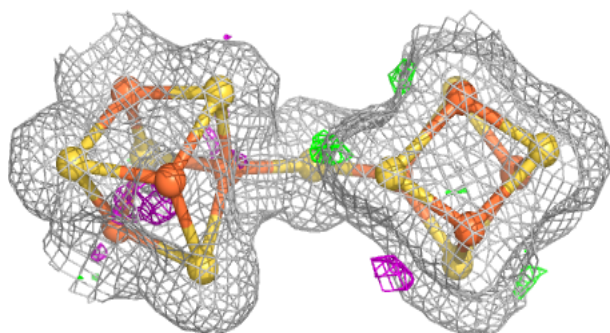
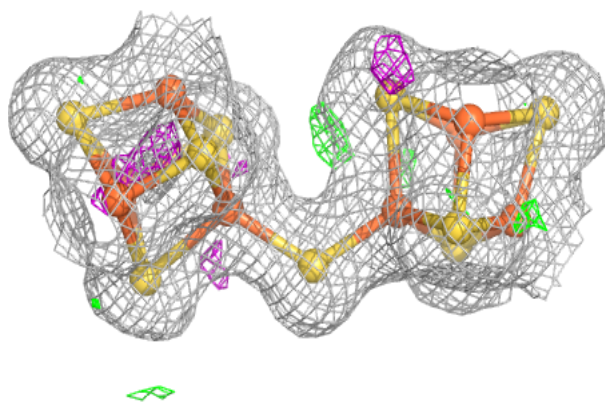
**Electron density around BJ8 C 501:**

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



Electron density around BJ8 D 501:

$2mF_o - DF_c$ (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.