



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:25 AM UTC

PDB ID : 6PAM / pdb_00006pam
Title : Structure of a bacterial Atm1-family ABC transporter with MgADP bound
Authors : Fan, C.; Kaiser, J.T.; Rees, D.C.
Deposited on : 2019-06-11
Resolution : 3.70 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

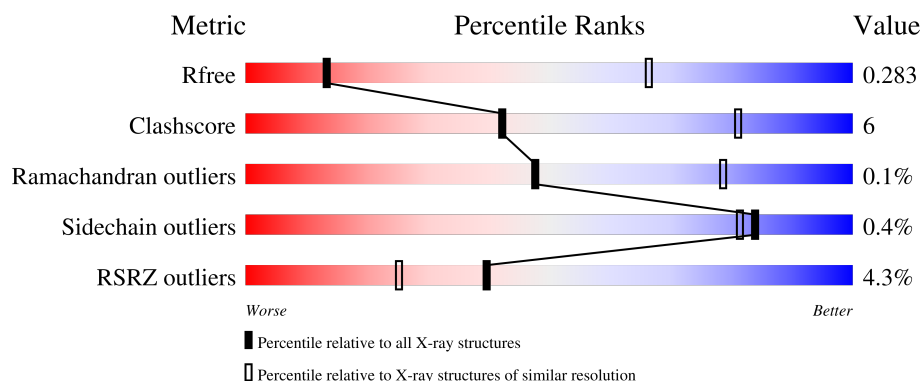
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 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	1131 (3.80-3.60)
Clashscore	190562	1171 (3.80-3.60)
Ramachandran outliers	187476	1129 (3.80-3.60)
Sidechain outliers	187428	1126 (3.80-3.60)
RSRZ outliers	180081	1130 (3.80-3.60)

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	614	3% 79% 16% 5%
1	B	614	3% 76% 19% 5%
1	C	614	4% 79% 17% .
1	D	614	4% 78% 17% 5%
1	E	614	4% 82% 14% .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	614	 6% 83% 13% •
1	G	614	 3% 82% 14% •
1	H	614	 6% 83% 13% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 36839 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 ATM1-type heavy metal exporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	586	Total	C	N	O	S	0	0	0
			4561	2913	812	823	13			
1	B	586	Total	C	N	O	S	0	0	0
			4561	2913	812	823	13			
1	C	590	Total	C	N	O	S	0	0	0
			4590	2929	816	832	13			
1	D	586	Total	C	N	O	S	0	0	0
			4561	2913	812	823	13			
1	E	589	Total	C	N	O	S	0	0	0
			4581	2924	815	829	13			
1	F	589	Total	C	N	O	S	0	0	0
			4581	2924	815	829	13			
1	G	590	Total	C	N	O	S	0	0	0
			4590	2929	816	832	13			
1	H	590	Total	C	N	O	S	0	0	0
			4590	2929	816	832	13			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	527	CYS	ALA	engineered mutation	UNP Q2G506
A	609	HIS	-	expression tag	UNP Q2G506
A	610	HIS	-	expression tag	UNP Q2G506
A	611	HIS	-	expression tag	UNP Q2G506
A	612	HIS	-	expression tag	UNP Q2G506
A	613	HIS	-	expression tag	UNP Q2G506
A	614	HIS	-	expression tag	UNP Q2G506
B	527	CYS	ALA	engineered mutation	UNP Q2G506
B	609	HIS	-	expression tag	UNP Q2G506
B	610	HIS	-	expression tag	UNP Q2G506
B	611	HIS	-	expression tag	UNP Q2G506
B	612	HIS	-	expression tag	UNP Q2G506
B	613	HIS	-	expression tag	UNP Q2G506

Continued on next page...

Continued from previous page...

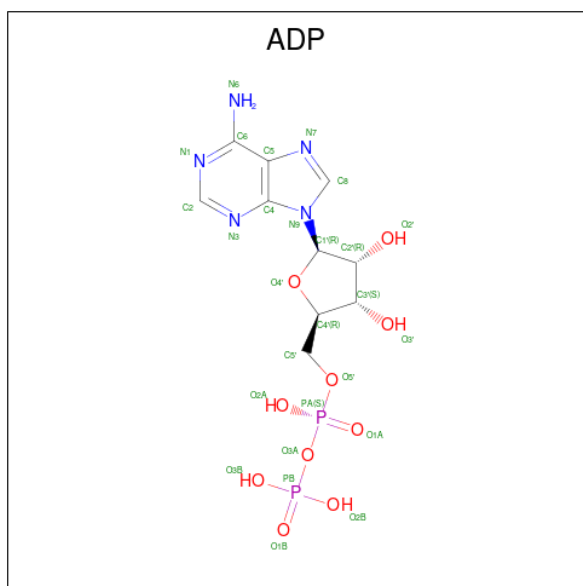
Chain	Residue	Modelled	Actual	Comment	Reference
B	614	HIS	-	expression tag	UNP Q2G506
C	527	CYS	ALA	engineered mutation	UNP Q2G506
C	609	HIS	-	expression tag	UNP Q2G506
C	610	HIS	-	expression tag	UNP Q2G506
C	611	HIS	-	expression tag	UNP Q2G506
C	612	HIS	-	expression tag	UNP Q2G506
C	613	HIS	-	expression tag	UNP Q2G506
C	614	HIS	-	expression tag	UNP Q2G506
D	527	CYS	ALA	engineered mutation	UNP Q2G506
D	609	HIS	-	expression tag	UNP Q2G506
D	610	HIS	-	expression tag	UNP Q2G506
D	611	HIS	-	expression tag	UNP Q2G506
D	612	HIS	-	expression tag	UNP Q2G506
D	613	HIS	-	expression tag	UNP Q2G506
D	614	HIS	-	expression tag	UNP Q2G506
E	527	CYS	ALA	engineered mutation	UNP Q2G506
E	609	HIS	-	expression tag	UNP Q2G506
E	610	HIS	-	expression tag	UNP Q2G506
E	611	HIS	-	expression tag	UNP Q2G506
E	612	HIS	-	expression tag	UNP Q2G506
E	613	HIS	-	expression tag	UNP Q2G506
E	614	HIS	-	expression tag	UNP Q2G506
F	527	CYS	ALA	engineered mutation	UNP Q2G506
F	609	HIS	-	expression tag	UNP Q2G506
F	610	HIS	-	expression tag	UNP Q2G506
F	611	HIS	-	expression tag	UNP Q2G506
F	612	HIS	-	expression tag	UNP Q2G506
F	613	HIS	-	expression tag	UNP Q2G506
F	614	HIS	-	expression tag	UNP Q2G506
G	527	CYS	ALA	engineered mutation	UNP Q2G506
G	609	HIS	-	expression tag	UNP Q2G506
G	610	HIS	-	expression tag	UNP Q2G506
G	611	HIS	-	expression tag	UNP Q2G506
G	612	HIS	-	expression tag	UNP Q2G506
G	613	HIS	-	expression tag	UNP Q2G506
G	614	HIS	-	expression tag	UNP Q2G506
H	527	CYS	ALA	engineered mutation	UNP Q2G506
H	609	HIS	-	expression tag	UNP Q2G506
H	610	HIS	-	expression tag	UNP Q2G506
H	611	HIS	-	expression tag	UNP Q2G506
H	612	HIS	-	expression tag	UNP Q2G506
H	613	HIS	-	expression tag	UNP Q2G506

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
H	614	HIS	-	expression tag	UNP Q2G506

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	G	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	H	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

Continued on next page...

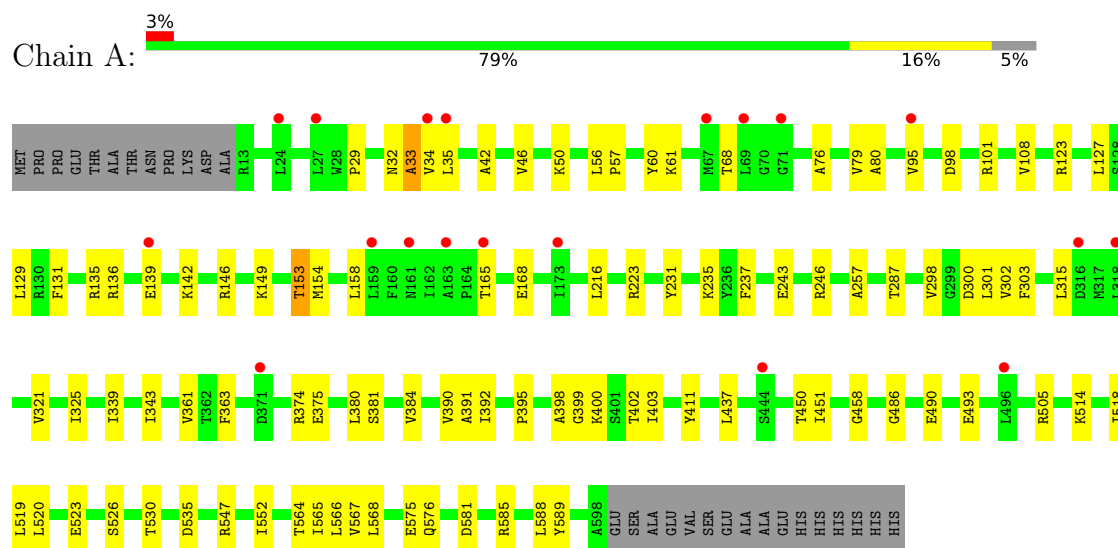
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Mg 1	0	0
3	C	1	Total 1	Mg 1	0	0
3	D	1	Total 1	Mg 1	0	0
3	E	1	Total 1	Mg 1	0	0
3	F	1	Total 1	Mg 1	0	0
3	G	1	Total 1	Mg 1	0	0
3	H	1	Total 1	Mg 1	0	0

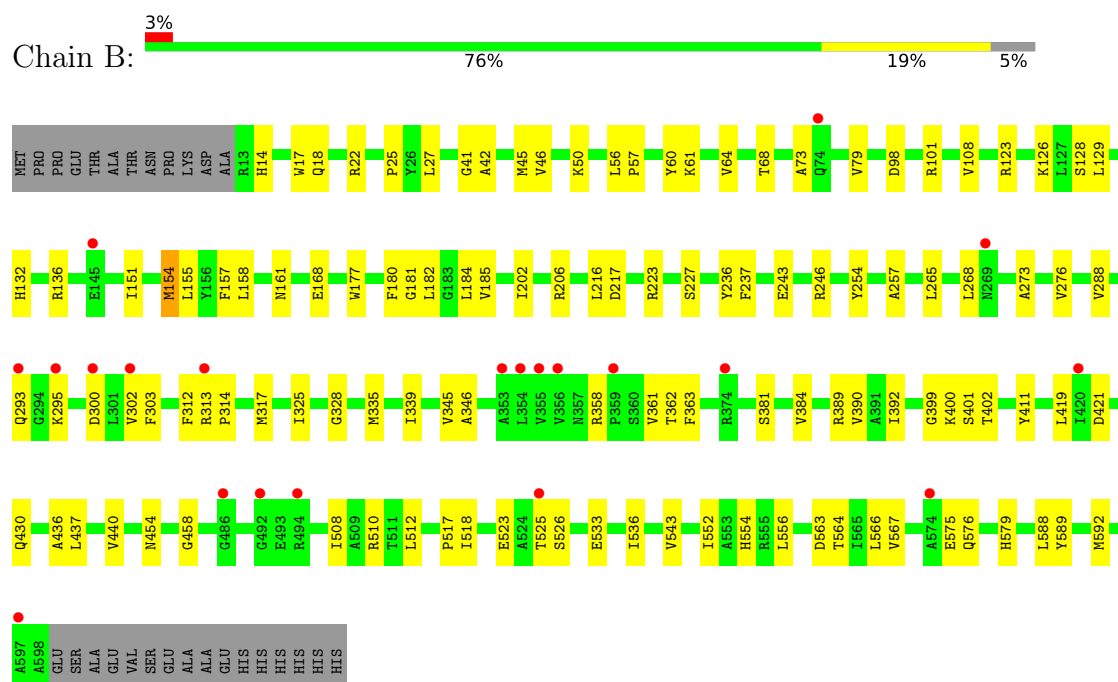
3 Residue-property plots [i](#)

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.

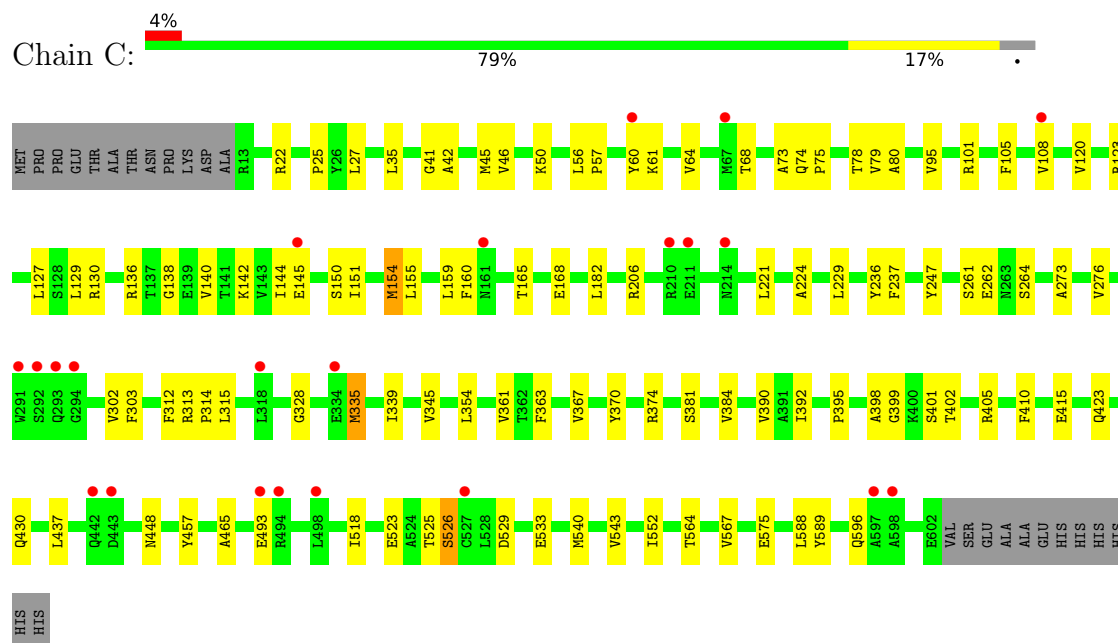
- Molecule 1: ATM1-type heavy metal exporter



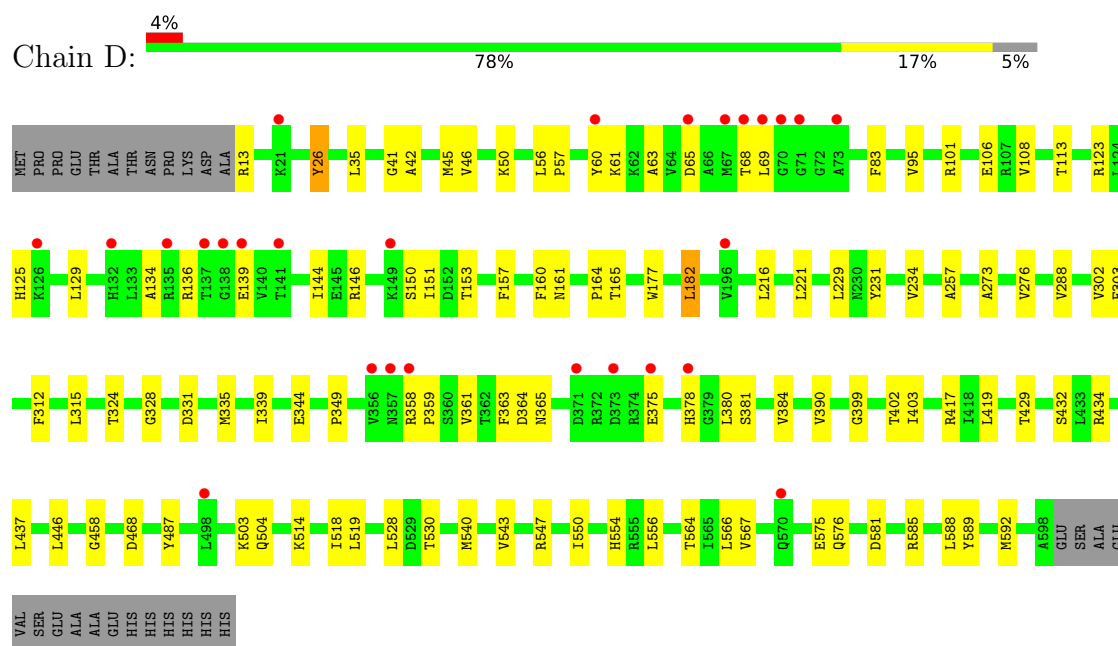
- Molecule 1: ATM1-type heavy metal exporter



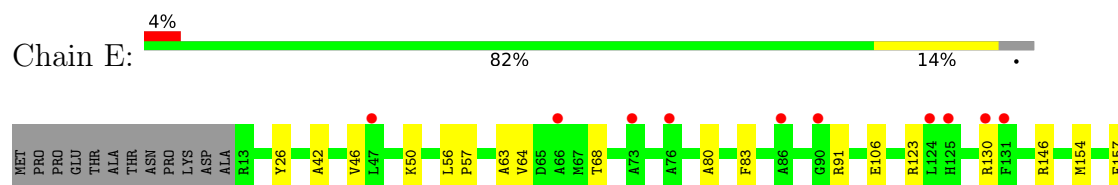
- Molecule 1: ATM1-type heavy metal exporter

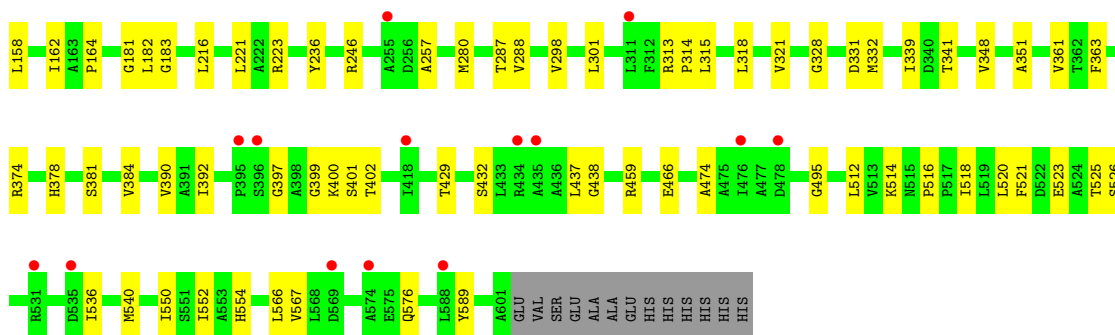


- Molecule 1: ATM1-type heavy metal exporter

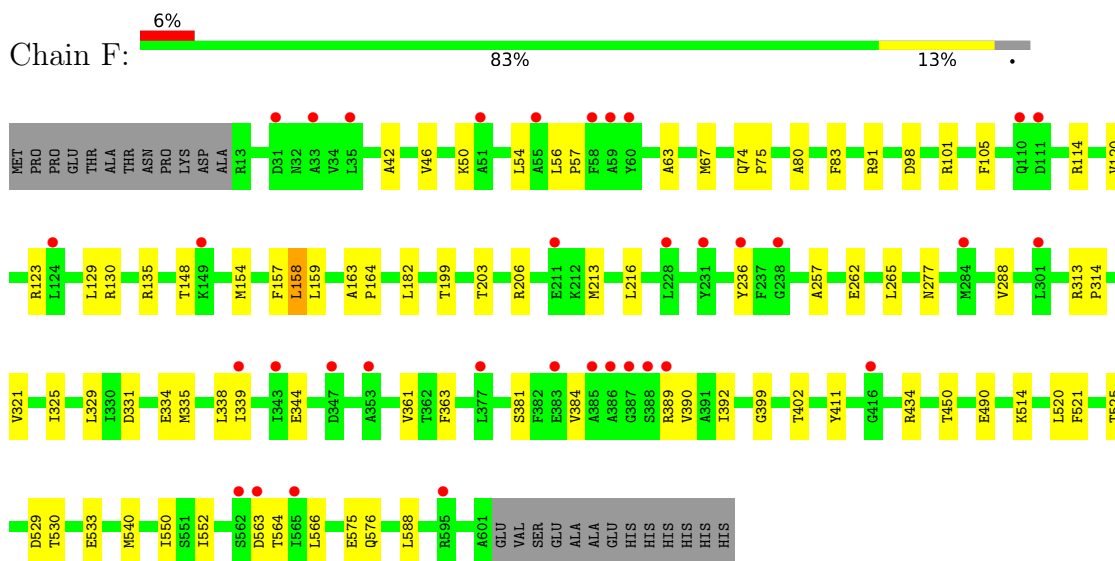


- Molecule 1: ATM1-type heavy metal exporter

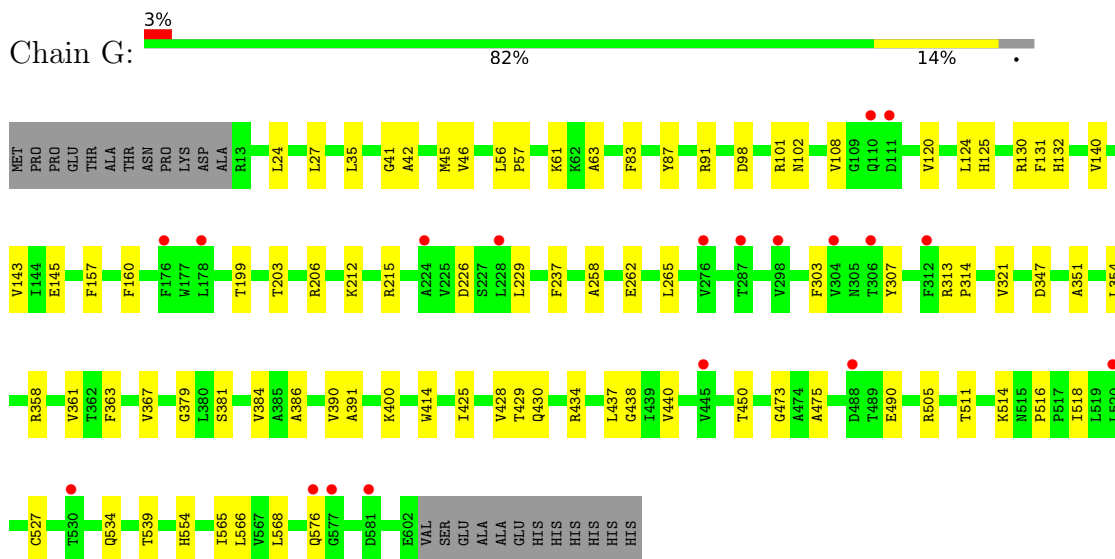




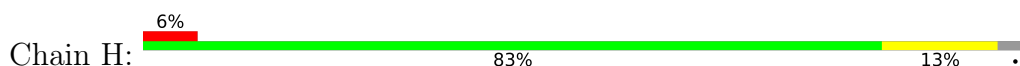
- Molecule 1: ATM1-type heavy metal exporter

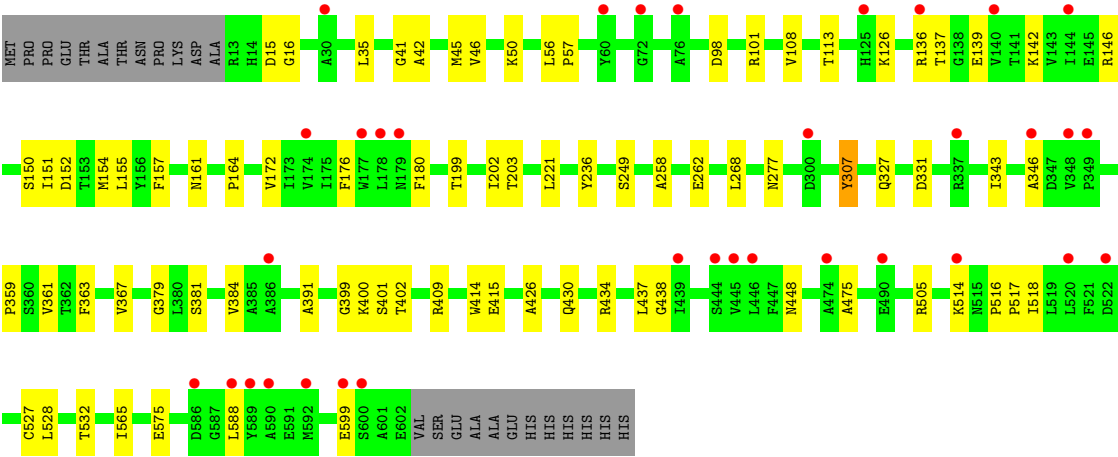


- Molecule 1: ATM1-type heavy metal exporter



- Molecule 1: ATM1-type heavy metal exporter





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	129.18Å 133.61Å 134.26Å 110.62° 98.28° 101.20°	Depositor
Resolution (Å)	39.61 – 3.70 39.61 – 3.70	Depositor EDS
% Data completeness (in resolution range)	97.6 (39.61-3.70) 97.7 (39.61-3.70)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 3.66Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472, REFMAC	Depositor
R, R_{free}	0.238 , 0.286 0.240 , 0.283	Depositor DCC
R_{free} test set	4195 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	146.7	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 83.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.055 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	36839	wwPDB-VP
Average B, all atoms (Å ²)	178.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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

5.1 Standard geometry

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

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	0.15	0/4646	0.36	2/6310 (0.0%)
1	B	0.15	0/4646	0.35	0/6310
1	C	0.14	0/4675	0.33	0/6349
1	D	0.13	0/4646	0.32	0/6310
1	E	0.13	0/4666	0.34	1/6337 (0.0%)
1	F	0.13	0/4666	0.32	0/6337
1	G	0.13	0/4675	0.33	0/6349
1	H	0.11	0/4675	0.29	0/6349
All	All	0.13	0/37295	0.33	3/50651 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	33	ALA	CA-C-N	5.20	131.32	121.97
1	A	33	ALA	C-N-CA	5.20	131.32	121.97
1	E	183	GLY	N-CA-C	-5.14	106.15	113.86

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	4561	0	4645	66	0
1	B	4561	0	4645	80	0
1	C	4590	0	4667	70	0
1	D	4561	0	4645	75	0
1	E	4581	0	4661	56	0
1	F	4581	0	4660	59	0
1	G	4590	0	4667	53	0
1	H	4590	0	4667	57	0
2	A	27	0	12	2	0
2	B	27	0	12	2	0
2	C	27	0	12	2	0
2	D	27	0	12	0	0
2	E	27	0	12	3	0
2	F	27	0	12	1	0
2	G	27	0	12	1	0
2	H	27	0	12	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
All	All	36839	0	37353	466	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (466) 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:G:354:LEU:HD11	1:G:428:VAL:HB	1.41	1.02
1:B:128:SER:HB3	1:B:345:VAL:HG21	1.59	0.84
1:D:146:ARG:NH1	1:D:335:MET:SD	2.55	0.79
1:F:154:MET:HG3	1:F:325:ILE:HA	1.65	0.79
1:A:136:ARG:HB2	1:A:139:GLU:HB2	1.64	0.78
1:C:73:ALA:HB1	1:C:79:VAL:HG21	1.65	0.78
1:B:525:THR:HG22	1:B:536:ILE:HD11	1.69	0.74
1:D:113:THR:HB	1:D:151:ILE:HD13	1.70	0.73
1:G:363:PHE:O	1:G:381:SER:HA	1.88	0.73
1:H:136:ARG:NH2	1:H:139:GLU:OE1	2.22	0.71

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:PHE:O	1:A:381:SER:HA	1.93	0.69
1:A:390:VAL:HG13	1:A:564:THR:HG23	1.74	0.68
1:A:29:PRO:HB2	1:A:32:ASN:HB3	1.75	0.68
1:A:35:LEU:HD12	1:A:35:LEU:H	1.57	0.68
1:C:374:ARG:HH22	2:C:701:ADP:H1'	1.59	0.67
1:D:567:VAL:HG21	1:D:589:TYR:HB2	1.75	0.67
1:E:363:PHE:O	1:E:381:SER:HA	1.95	0.67
1:E:315:LEU:HD22	1:E:318:LEU:HD22	1.75	0.67
1:E:123:ARG:NH2	1:E:341:THR:O	2.28	0.67
1:F:154:MET:HE2	1:F:158:LEU:HD12	1.75	0.67
1:E:123:ARG:HG2	1:E:339:ILE:HA	1.77	0.66
1:B:177:TRP:HA	1:B:181:GLY:HA3	1.76	0.66
1:D:153:THR:HB	1:D:324:THR:HB	1.76	0.66
1:E:236:TYR:O	1:F:514:LYS:NZ	2.28	0.65
1:G:527:CYS:HB3	1:H:527:CYS:HA	1.79	0.65
1:B:18:GLN:NE2	1:D:364:ASP:OD1	2.28	0.65
1:C:229:LEU:HD11	1:D:136:ARG:HH22	1.62	0.65
1:F:123:ARG:HG3	1:F:339:ILE:HG23	1.78	0.65
1:A:149:LYS:O	1:A:153:THR:HG23	1.97	0.64
1:G:226:ASP:OD2	1:H:448:ASN:N	2.25	0.64
1:E:567:VAL:HG21	1:E:589:TYR:HB2	1.80	0.64
1:D:177:TRP:HZ3	1:D:182:LEU:HD13	1.63	0.64
1:A:68:THR:HG21	1:A:298:VAL:HG11	1.81	0.63
1:A:165:THR:HG21	1:A:315:LEU:HD21	1.80	0.63
1:D:384:VAL:HG11	1:D:550:ILE:HD11	1.79	0.63
1:F:216:LEU:HD12	1:F:257:ALA:HB2	1.80	0.63
1:G:46:VAL:HG22	1:G:101:ARG:HD2	1.79	0.63
1:E:401:SER:OG	2:E:701:ADP:O1A	2.16	0.63
1:B:400:LYS:NZ	2:B:701:ADP:O1B	2.30	0.63
1:G:91:ARG:NH1	1:H:277:ASN:OD1	2.33	0.62
1:E:392:ILE:HB	1:E:552:ILE:HG12	1.80	0.62
1:E:64:VAL:O	1:E:68:THR:HG23	1.99	0.62
1:B:128:SER:O	1:B:129:LEU:HG	1.98	0.62
1:C:236:TYR:O	1:D:514:LYS:NZ	2.29	0.62
1:B:123:ARG:HG2	1:B:339:ILE:HA	1.82	0.62
1:C:523:GLU:HB3	1:C:526:SER:OG	1.98	0.61
1:C:363:PHE:O	1:C:381:SER:HA	2.01	0.61
1:C:75:PRO:HG2	1:C:78:THR:HB	1.83	0.61
1:E:525:THR:HG22	1:E:536:ILE:HD11	1.83	0.61
1:C:567:VAL:HG21	1:C:589:TYR:HB2	1.83	0.60
1:E:429:THR:HG1	1:E:432:SER:HG	1.49	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:ARG:NH2	2:A:701:ADP:N3	2.49	0.60
1:C:575:GLU:OE2	1:C:588:LEU:N	2.33	0.60
1:H:146:ARG:O	1:H:150:SER:N	2.34	0.60
1:A:154:MET:HG3	1:A:325:ILE:HA	1.83	0.60
1:D:150:SER:HB2	1:D:331:ASP:HB2	1.83	0.60
1:G:514:LYS:NZ	1:H:236:TYR:O	2.34	0.60
1:A:123:ARG:HG2	1:A:339:ILE:HA	1.83	0.60
1:F:361:VAL:HB	1:F:384:VAL:HB	1.83	0.60
1:A:129:LEU:HD13	1:A:411:TYR:HE2	1.66	0.60
1:F:206:ARG:HD2	1:F:265:LEU:HA	1.83	0.60
1:G:63:ALA:HB2	1:G:83:PHE:HB3	1.83	0.60
1:G:56:LEU:HG	1:G:57:PRO:HD3	1.84	0.59
1:C:448:ASN:ND2	1:C:493:GLU:OE2	2.35	0.59
1:E:91:ARG:NH1	1:F:277:ASN:OD1	2.35	0.59
1:B:64:VAL:O	1:B:68:THR:HG23	2.02	0.59
1:D:399:GLY:O	1:D:402:THR:OG1	2.15	0.59
1:B:361:VAL:HB	1:B:384:VAL:HB	1.84	0.58
1:D:504:GLN:HG3	1:D:528:LEU:HD11	1.84	0.58
1:E:554:HIS:CE1	1:F:530:THR:H	2.22	0.57
1:B:567:VAL:HG21	1:B:589:TYR:HB2	1.85	0.57
1:H:98:ASP:OD1	1:H:101:ARG:NH1	2.38	0.57
1:C:56:LEU:HG	1:C:57:PRO:HD3	1.87	0.57
1:B:73:ALA:HB2	1:B:79:VAL:HG23	1.85	0.57
1:D:177:TRP:CZ3	1:D:182:LEU:HD13	2.40	0.57
1:E:280:MET:HE3	1:F:91:ARG:HH12	1.69	0.57
1:C:127:LEU:O	1:D:231:TYR:OH	2.19	0.57
1:E:374:ARG:NH1	2:E:701:ADP:O3'	2.38	0.57
1:C:101:ARG:HE	1:C:160:PHE:HD2	1.51	0.56
1:D:56:LEU:HG	1:D:57:PRO:HD3	1.87	0.56
1:H:391:ALA:HB3	1:H:565:ILE:HG12	1.87	0.56
1:D:123:ARG:HG3	1:D:339:ILE:HG23	1.88	0.56
1:F:363:PHE:O	1:F:381:SER:HA	2.06	0.56
1:F:399:GLY:O	1:F:402:THR:OG1	2.22	0.56
1:G:157:PHE:HD1	1:G:321:VAL:HA	1.70	0.56
1:A:80:ALA:HB1	1:B:288:VAL:HG21	1.88	0.56
1:A:399:GLY:O	1:A:402:THR:OG1	2.20	0.56
1:C:361:VAL:HB	1:C:384:VAL:HB	1.87	0.56
1:E:50:LYS:HE3	1:E:164:PRO:HB3	1.86	0.56
1:F:450:THR:HA	1:F:490:GLU:HA	1.87	0.56
1:D:101:ARG:HE	1:D:160:PHE:HD1	1.52	0.56
1:B:401:SER:OG	2:B:701:ADP:O1A	2.21	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:PHE:HE2	1:B:182:LEU:HB2	1.71	0.56
1:F:50:LYS:HE3	1:F:164:PRO:HB3	1.88	0.56
1:H:475:ALA:O	1:H:505:ARG:NH1	2.38	0.56
1:F:46:VAL:HG22	1:F:101:ARG:HD2	1.88	0.55
1:H:56:LEU:HG	1:H:57:PRO:HD3	1.88	0.55
1:B:50:LYS:NZ	1:B:168:GLU:OE1	2.37	0.55
1:G:361:VAL:HB	1:G:384:VAL:HB	1.87	0.55
1:C:64:VAL:O	1:C:68:THR:HG23	2.06	0.55
1:G:554:HIS:NE2	2:G:701:ADP:O2B	2.33	0.55
1:E:361:VAL:HB	1:E:384:VAL:HB	1.89	0.55
1:C:392:ILE:HB	1:C:552:ILE:HG12	1.88	0.55
1:E:521:PHE:CD2	1:E:540:MET:HE3	2.42	0.55
1:E:56:LEU:HG	1:E:57:PRO:HD3	1.89	0.54
1:B:56:LEU:HG	1:B:57:PRO:HD3	1.89	0.54
1:B:525:THR:HB	1:B:533:GLU:HG3	1.89	0.54
1:F:63:ALA:HB2	1:F:83:PHE:HB3	1.88	0.54
1:H:363:PHE:O	1:H:381:SER:HA	2.07	0.54
1:C:151:ILE:HD12	1:C:335:MET:HE3	1.88	0.54
1:C:27:LEU:HD21	1:C:155:LEU:HD21	1.88	0.54
1:G:475:ALA:O	1:G:505:ARG:NH1	2.41	0.54
1:H:41:GLY:O	1:H:45:MET:HG2	2.08	0.54
1:H:50:LYS:HE3	1:H:164:PRO:HB3	1.90	0.54
1:B:358:ARG:N	1:B:421:ASP:OD1	2.37	0.54
1:F:157:PHE:HD2	1:F:321:VAL:HG22	1.71	0.54
1:A:361:VAL:HB	1:A:384:VAL:HB	1.90	0.54
1:B:18:GLN:HE22	1:D:365:ASN:HB2	1.73	0.54
1:F:154:MET:HE3	1:F:325:ILE:HG23	1.90	0.54
1:B:399:GLY:O	1:B:402:THR:OG1	2.21	0.54
1:G:425:ILE:O	1:G:428:VAL:HG22	2.08	0.54
1:A:56:LEU:HG	1:A:57:PRO:HD3	1.90	0.54
1:E:315:LEU:HA	1:E:318:LEU:HB2	1.88	0.53
1:A:575:GLU:OE2	1:A:588:LEU:N	2.37	0.53
1:B:384:VAL:HG13	1:B:390:VAL:HG21	1.91	0.53
1:F:56:LEU:HG	1:F:57:PRO:HD3	1.90	0.53
1:G:450:THR:HA	1:G:490:GLU:HA	1.90	0.53
1:C:221:LEU:HD21	1:D:144:ILE:HA	1.90	0.53
1:F:129:LEU:HD12	1:F:411:TYR:HE2	1.73	0.53
1:H:172:VAL:HG22	1:H:307:TYR:HD2	1.74	0.53
1:C:140:VAL:HG22	1:D:229:LEU:HD21	1.91	0.53
1:D:575:GLU:OE2	1:D:588:LEU:N	2.40	0.53
1:H:146:ARG:NH1	1:H:331:ASP:OD2	2.42	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LYS:HB2	1:A:303:PHE:HB2	1.90	0.53
1:H:202:ILE:HG21	1:H:268:LEU:HB2	1.91	0.53
1:H:151:ILE:HA	1:H:154:MET:HG2	1.91	0.53
1:A:33:ALA:O	1:A:35:LEU:N	2.42	0.52
1:A:142:LYS:O	1:A:146:ARG:HB3	2.09	0.52
1:G:354:LEU:CD1	1:G:428:VAL:HB	2.25	0.52
1:H:35:LEU:HD22	1:H:108:VAL:HG22	1.91	0.52
1:B:158:LEU:HD23	1:B:325:ILE:HD13	1.91	0.52
1:A:451:ILE:HD11	1:A:486:GLY:O	2.09	0.52
1:B:512:LEU:HD21	1:B:543:VAL:HG13	1.91	0.52
1:H:367:VAL:HG22	1:H:379:GLY:H	1.74	0.52
1:F:392:ILE:HB	1:F:552:ILE:HG12	1.92	0.52
1:C:395:PRO:HD2	1:C:398:ALA:HB2	1.92	0.52
1:B:392:ILE:HB	1:B:552:ILE:HG12	1.92	0.52
1:H:505:ARG:NH2	1:H:532:THR:OG1	2.43	0.52
1:D:361:VAL:HB	1:D:384:VAL:HB	1.91	0.52
1:B:129:LEU:HA	1:B:132:HIS:ND1	2.24	0.52
1:G:199:THR:O	1:G:203:THR:OG1	2.20	0.52
1:A:223:ARG:HD3	1:A:246:ARG:HG2	1.92	0.52
1:B:14:HIS:NE2	1:B:17:TRP:HB2	2.25	0.51
1:H:142:LYS:HE3	1:H:146:ARG:HH21	1.75	0.51
1:H:438:GLY:HA3	1:H:516:PRO:HG3	1.91	0.51
1:B:363:PHE:O	1:B:381:SER:HA	2.10	0.51
1:F:74:GLN:HB2	1:F:75:PRO:HD3	1.93	0.51
1:C:402:THR:HA	1:C:405:ARG:HG2	1.91	0.51
1:A:567:VAL:HG21	1:A:589:TYR:HB2	1.92	0.51
1:H:361:VAL:HB	1:H:384:VAL:HB	1.93	0.51
1:H:157:PHE:O	1:H:161:ASN:HB3	2.11	0.51
1:B:126:LYS:NZ	1:D:349:PRO:O	2.44	0.51
1:B:129:LEU:HD22	1:B:411:TYR:CE2	2.45	0.51
1:G:140:VAL:HA	1:G:143:VAL:HG22	1.92	0.51
1:C:74:GLN:HB2	1:C:75:PRO:HD3	1.92	0.51
1:D:363:PHE:O	1:D:381:SER:HA	2.11	0.51
1:B:575:GLU:OE2	1:B:588:LEU:N	2.42	0.51
1:E:474:ALA:HB2	1:E:512:LEU:HD12	1.93	0.51
1:A:375:GLU:OE1	1:E:378:HIS:NE2	2.44	0.50
1:A:520:LEU:HD11	1:A:552:ILE:HD12	1.91	0.50
1:C:35:LEU:HD22	1:C:108:VAL:HG12	1.93	0.50
1:C:276:VAL:HG13	1:C:312:PHE:HD2	1.75	0.50
1:E:523:GLU:HB3	1:E:526:SER:OG	2.11	0.50
1:C:384:VAL:HG13	1:C:390:VAL:HG21	1.91	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:384:VAL:HG13	1:E:390:VAL:HG21	1.93	0.50
1:A:380:LEU:HD23	1:A:403:ILE:HD11	1.93	0.50
1:H:367:VAL:O	1:H:414:TRP:N	2.44	0.50
1:A:523:GLU:HB3	1:A:526:SER:OG	2.12	0.50
1:G:438:GLY:HA3	1:G:516:PRO:HG3	1.93	0.50
1:A:530:THR:H	1:B:554:HIS:CE1	2.30	0.50
1:D:434:ARG:O	1:D:514:LYS:NZ	2.43	0.50
1:A:60:TYR:CE1	1:A:302:VAL:HG11	2.47	0.49
1:C:151:ILE:CD1	1:C:335:MET:HE3	2.43	0.49
1:D:429:THR:HG1	1:D:432:SER:HG	1.59	0.49
1:E:288:VAL:HG21	1:F:80:ALA:HB1	1.93	0.49
1:B:180:PHE:CE2	1:B:182:LEU:HB2	2.48	0.49
1:C:165:THR:HG21	1:C:315:LEU:HD21	1.94	0.49
1:F:199:THR:O	1:F:203:THR:OG1	2.27	0.49
1:B:206:ARG:HD2	1:B:265:LEU:HA	1.93	0.49
1:C:136:ARG:O	1:C:140:VAL:HG23	2.12	0.49
1:D:46:VAL:HG22	1:D:101:ARG:HD2	1.95	0.49
1:G:347:ASP:HB3	1:G:430:GLN:HG3	1.94	0.49
1:B:440:VAL:HG21	1:B:508:ILE:HD13	1.94	0.49
1:F:331:ASP:O	1:F:335:MET:HG2	2.13	0.49
1:B:362:THR:HB	1:B:419:LEU:HB2	1.94	0.49
1:D:165:THR:HG21	1:D:315:LEU:HD21	1.95	0.49
1:F:154:MET:CE	1:F:329:LEU:HG	2.43	0.49
1:A:50:LYS:NZ	1:A:168:GLU:OE1	2.44	0.49
1:B:151:ILE:HD11	1:B:335:MET:SD	2.53	0.49
1:B:566:LEU:HD23	1:B:576:GLN:HG3	1.95	0.49
1:D:358:ARG:HD2	1:D:359:PRO:HD2	1.94	0.49
1:A:400:LYS:HB3	1:A:568:LEU:HD12	1.95	0.48
1:D:63:ALA:HB2	1:D:83:PHE:HB3	1.94	0.48
1:F:434:ARG:O	1:F:514:LYS:NZ	2.42	0.48
1:F:521:PHE:CD2	1:F:540:MET:HE3	2.48	0.48
1:E:130:ARG:HH21	1:E:341:THR:HG21	1.78	0.48
1:A:395:PRO:HD2	1:A:398:ALA:HB2	1.95	0.48
1:C:144:ILE:HG13	1:C:145:GLU:H	1.79	0.48
1:F:130:ARG:NH1	1:F:338:LEU:HD22	2.29	0.48
1:B:129:LEU:HA	1:B:132:HIS:H	1.79	0.48
1:C:144:ILE:HB	1:D:221:LEU:HG	1.94	0.48
1:G:367:VAL:HG22	1:G:379:GLY:H	1.78	0.48
1:C:150:SER:OG	1:C:328:GLY:HA2	2.13	0.48
1:G:440:VAL:HG22	1:G:511:THR:HG21	1.95	0.48
1:H:172:VAL:HG22	1:H:307:TYR:CD2	2.48	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:298:VAL:HG23	1:F:67:MET:HE2	1.95	0.48
1:G:367:VAL:O	1:G:414:TRP:N	2.45	0.48
1:D:65:ASP:HA	1:D:68:THR:HG22	1.94	0.48
1:E:459:ARG:NH2	1:E:466:GLU:OE2	2.35	0.48
1:A:391:ALA:HB3	1:A:565:ILE:HD13	1.94	0.48
1:A:392:ILE:HB	1:A:552:ILE:HG12	1.96	0.47
1:C:80:ALA:HB1	1:D:288:VAL:HG21	1.96	0.47
1:D:68:THR:HG23	1:D:69:LEU:HD13	1.95	0.47
1:A:321:VAL:O	1:A:325:ILE:HG12	2.14	0.47
1:D:113:THR:HG21	1:D:151:ILE:HG21	1.95	0.47
1:E:63:ALA:HB2	1:E:83:PHE:HB3	1.96	0.47
1:E:399:GLY:O	1:E:402:THR:OG1	2.29	0.47
1:C:60:TYR:CE1	1:C:302:VAL:HG11	2.48	0.47
1:H:346:ALA:O	1:H:409:ARG:NH2	2.47	0.47
1:C:465:ALA:H	1:H:249:SER:HB2	1.80	0.47
1:D:390:VAL:HG13	1:D:564:THR:HG23	1.97	0.47
1:E:80:ALA:HB1	1:F:288:VAL:HG21	1.95	0.47
1:G:42:ALA:O	1:G:46:VAL:HG23	2.14	0.47
1:F:157:PHE:CD2	1:F:321:VAL:HG22	2.49	0.47
1:B:129:LEU:HB2	1:B:132:HIS:HB2	1.96	0.47
1:E:400:LYS:NZ	1:E:523:GLU:OE1	2.47	0.47
1:F:154:MET:O	1:F:158:LEU:HB2	2.15	0.47
1:G:384:VAL:HG13	1:G:390:VAL:HG11	1.95	0.47
1:A:231:TYR:OH	1:A:235:LYS:HE2	2.14	0.47
1:B:276:VAL:HG13	1:B:312:PHE:HD2	1.80	0.47
1:C:367:VAL:HB	1:C:415:GLU:HB2	1.96	0.47
1:C:399:GLY:O	1:C:402:THR:OG1	2.22	0.47
1:F:399:GLY:HA2	2:F:701:ADP:H5'2	1.95	0.47
1:G:430:GLN:HE21	1:G:434:ARG:HH22	1.62	0.47
1:A:95:VAL:HG11	1:B:273:ALA:HB2	1.97	0.47
1:A:505:ARG:HH22	1:A:535:ASP:HB3	1.79	0.47
1:A:519:LEU:HD13	1:A:547:ARG:HD3	1.96	0.47
1:F:575:GLU:OE2	1:F:588:LEU:N	2.44	0.47
1:G:87:TYR:CZ	1:G:91:ARG:HD2	2.50	0.47
1:D:113:THR:HG21	1:D:151:ILE:CG2	2.45	0.47
1:D:216:LEU:HD12	1:D:257:ALA:HB2	1.97	0.47
1:A:437:LEU:HD23	1:A:518:ILE:HB	1.96	0.46
1:G:132:HIS:O	1:G:132:HIS:ND1	2.48	0.46
1:H:45:MET:HB2	1:H:101:ARG:HB3	1.97	0.46
1:E:566:LEU:HD23	1:E:576:GLN:HG3	1.97	0.46
1:H:437:LEU:HD23	1:H:518:ILE:HB	1.96	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:VAL:HG13	1:A:390:VAL:HG21	1.97	0.46
1:B:216:LEU:HD12	1:B:257:ALA:HB2	1.97	0.46
1:C:41:GLY:O	1:C:45:MET:HG2	2.15	0.46
1:D:384:VAL:HG13	1:D:390:VAL:HG21	1.97	0.46
1:A:32:ASN:ND2	1:A:35:LEU:HD13	2.30	0.46
1:A:98:ASP:O	1:A:101:ARG:HG2	2.15	0.46
1:B:227:SER:OG	1:B:243:GLU:HG2	2.16	0.46
1:C:50:LYS:NZ	1:C:168:GLU:OE1	2.43	0.46
1:H:126:LYS:HD3	1:H:343:ILE:HD11	1.97	0.46
1:B:293:GLN:OE1	1:B:295:LYS:NZ	2.48	0.46
1:C:370:TYR:OH	1:C:405:ARG:NH1	2.48	0.46
1:F:42:ALA:O	1:F:46:VAL:HG23	2.15	0.46
1:A:46:VAL:HG22	1:A:101:ARG:HD2	1.97	0.46
1:A:392:ILE:HG12	1:A:566:LEU:HD12	1.98	0.46
1:H:146:ARG:NH1	1:H:327:GLN:OE1	2.49	0.46
1:H:359:PRO:HB2	1:H:518:ILE:HD11	1.97	0.46
1:B:27:LEU:HD21	1:B:155:LEU:HD21	1.98	0.46
1:B:154:MET:HB2	1:B:328:GLY:HA3	1.97	0.46
1:C:237:PHE:CD1	1:D:458:GLY:HA2	2.51	0.46
1:E:348:VAL:HG23	1:E:351:ALA:HB2	1.98	0.46
1:F:120:VAL:HG11	1:F:335:MET:HE3	1.96	0.46
1:G:35:LEU:HD22	1:G:108:VAL:HG22	1.97	0.46
1:G:145:GLU:CD	1:H:221:LEU:HD21	2.41	0.46
1:D:276:VAL:HG13	1:D:312:PHE:HD2	1.81	0.46
1:G:212:LYS:HG3	1:G:215:ARG:NH2	2.30	0.46
1:B:98:ASP:O	1:B:101:ARG:HG2	2.16	0.45
1:C:27:LEU:HD11	1:C:155:LEU:HD11	1.97	0.45
1:B:436:ALA:HA	1:B:517:PRO:HD2	1.97	0.45
1:B:556:LEU:HD11	1:B:592:MET:HB3	1.98	0.45
1:D:380:LEU:HD23	1:D:403:ILE:HD11	1.97	0.45
1:H:154:MET:HG3	1:H:155:LEU:N	2.31	0.45
1:H:528:LEU:HD13	1:H:532:THR:HG22	1.98	0.45
1:C:61:LYS:HB2	1:C:303:PHE:HB2	1.98	0.45
1:C:390:VAL:HG13	1:C:564:THR:HG23	1.98	0.45
1:D:556:LEU:HD11	1:D:592:MET:HB3	1.97	0.45
1:E:158:LEU:HD23	1:E:162:ILE:HG13	1.97	0.45
1:F:344:GLU:O	1:F:344:GLU:HG2	2.16	0.45
1:G:400:LYS:HB3	1:G:568:LEU:HD12	1.98	0.45
1:A:61:LYS:NZ	1:A:300:ASP:OD1	2.38	0.45
1:C:525:THR:HB	1:C:533:GLU:HG3	1.97	0.45
1:C:46:VAL:HG22	1:C:101:ARG:HD2	1.99	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154:MET:HB2	1:C:328:GLY:HA3	1.99	0.45
1:B:437:LEU:HD23	1:B:518:ILE:HB	1.97	0.45
1:B:523:GLU:HB3	1:B:526:SER:HB2	1.99	0.45
1:G:358:ARG:HG3	1:G:386:ALA:HB1	1.98	0.45
1:G:61:LYS:HB2	1:G:303:PHE:HB2	1.99	0.45
1:H:399:GLY:O	1:H:402:THR:OG1	2.32	0.45
1:C:130:ARG:HA	1:C:130:ARG:HD2	1.83	0.45
1:C:273:ALA:HB2	1:D:95:VAL:HG11	1.98	0.45
1:F:390:VAL:HG13	1:F:564:THR:HG23	1.98	0.45
1:H:359:PRO:HG3	1:H:517:PRO:HB2	1.99	0.45
1:D:129:LEU:HD13	1:D:344:GLU:OE2	2.17	0.44
1:F:120:VAL:HG11	1:F:335:MET:CE	2.47	0.44
1:A:514:LYS:NZ	1:B:236:TYR:O	2.45	0.44
1:B:22:ARG:HH22	1:D:419:LEU:HD11	1.82	0.44
1:D:50:LYS:HE3	1:D:164:PRO:HB3	1.98	0.44
1:E:438:GLY:HA3	1:E:516:PRO:HG3	2.00	0.44
1:H:142:LYS:HG2	1:H:146:ARG:HH21	1.82	0.44
1:H:199:THR:O	1:H:203:THR:OG1	2.32	0.44
1:A:127:LEU:HD23	1:A:343:ILE:HG12	1.99	0.44
1:D:157:PHE:O	1:D:161:ASN:HB3	2.18	0.44
1:G:41:GLY:O	1:G:45:MET:HG2	2.18	0.44
1:H:409:ARG:NE	1:H:426:ALA:HA	2.31	0.44
1:D:61:LYS:HB2	1:D:303:PHE:HB2	2.00	0.44
1:F:313:ARG:HB3	1:F:314:PRO:HD3	2.00	0.44
1:B:60:TYR:CE1	1:B:302:VAL:HG11	2.52	0.44
1:C:22:ARG:O	1:C:25:PRO:HD2	2.17	0.44
1:F:384:VAL:HG13	1:F:390:VAL:HG21	1.98	0.44
1:H:575:GLU:OE2	1:H:588:LEU:N	2.42	0.44
1:A:131:PHE:O	1:A:135:ARG:HG2	2.17	0.44
1:D:468:ASP:OD1	1:D:487:TYR:OH	2.30	0.44
1:F:105:PHE:CZ	1:F:159:LEU:HD12	2.53	0.44
1:G:102:ASN:OD1	1:G:160:PHE:HE2	2.01	0.44
1:G:351:ALA:HB3	1:G:429:THR:HG23	1.99	0.44
1:H:176:PHE:HA	1:H:180:PHE:HD2	1.82	0.44
1:C:123:ARG:HG2	1:C:339:ILE:HA	1.99	0.44
1:B:61:LYS:NZ	1:B:300:ASP:OD1	2.35	0.44
1:D:375:GLU:HG2	1:D:378:HIS:HD2	1.83	0.44
1:G:98:ASP:O	1:G:101:ARG:HG2	2.18	0.44
1:H:528:LEU:HD22	1:H:532:THR:HG21	2.00	0.44
1:E:313:ARG:HB3	1:E:314:PRO:HD3	1.99	0.43
1:G:258:ALA:O	1:G:262:GLU:HG2	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:GLY:O	1:B:45:MET:HG2	2.18	0.43
1:B:389:ARG:O	1:B:563:ASP:N	2.51	0.43
1:C:345:VAL:O	1:C:430:GLN:NE2	2.47	0.43
1:E:157:PHE:CD1	1:E:321:VAL:HG22	2.53	0.43
1:E:288:VAL:HG23	1:F:67:MET:HE3	2.00	0.43
1:G:212:LYS:HG3	1:G:215:ARG:HH22	1.82	0.43
1:A:374:ARG:HH22	2:A:701:ADP:HI'	1.84	0.43
1:C:120:VAL:HG21	1:C:335:MET:SD	2.58	0.43
1:E:437:LEU:HD23	1:E:518:ILE:HB	1.99	0.43
1:B:22:ARG:O	1:B:25:PRO:HD2	2.18	0.43
1:C:105:PHE:CZ	1:C:159:LEU:HD12	2.53	0.43
1:F:334:GLU:O	1:F:338:LEU:HG	2.19	0.43
1:C:529:ASP:HA	1:D:554:HIS:CE1	2.54	0.43
1:F:163:ALA:HB3	1:F:164:PRO:HD3	2.01	0.43
1:A:237:PHE:CD1	1:B:458:GLY:HA2	2.54	0.43
1:B:390:VAL:HG13	1:B:564:THR:HG23	1.99	0.43
1:E:154:MET:HB2	1:E:328:GLY:HA3	2.00	0.43
1:H:367:VAL:HB	1:H:415:GLU:HB2	2.00	0.43
1:D:566:LEU:HD23	1:D:576:GLN:HG3	1.99	0.43
1:D:26:TYR:N	1:D:26:TYR:CD2	2.87	0.43
1:D:358:ARG:HD2	1:D:358:ARG:HA	1.86	0.43
1:E:146:ARG:NH2	1:E:331:ASP:OD2	2.52	0.43
1:F:389:ARG:O	1:F:563:ASP:N	2.52	0.43
1:H:258:ALA:O	1:H:262:GLU:HG2	2.19	0.43
1:B:202:ILE:HB	1:B:268:LEU:HD13	2.00	0.42
1:D:134:ALA:HB1	1:D:139:GLU:CD	2.44	0.42
1:E:495:GLY:O	1:F:135:ARG:NH2	2.46	0.42
1:A:566:LEU:HD23	1:A:576:GLN:HG3	2.01	0.42
1:B:42:ALA:O	1:B:46:VAL:HG23	2.19	0.42
1:B:206:ARG:NH1	1:B:265:LEU:HB2	2.33	0.42
1:D:42:ALA:O	1:D:46:VAL:HG23	2.20	0.42
1:D:519:LEU:HD13	1:D:547:ARG:HD3	2.01	0.42
1:E:42:ALA:O	1:E:46:VAL:HG23	2.18	0.42
1:A:523:GLU:HB3	1:A:526:SER:HG	1.84	0.42
1:C:206:ARG:HG3	1:C:264:SER:OG	2.20	0.42
1:C:313:ARG:HB3	1:C:314:PRO:HD3	2.01	0.42
1:E:520:LEU:HD11	1:E:550:ILE:HD12	2.01	0.42
1:F:154:MET:HE3	1:F:329:LEU:HG	2.02	0.42
1:F:525:THR:HB	1:F:533:GLU:HG3	2.01	0.42
1:C:42:ALA:O	1:C:46:VAL:HG23	2.20	0.42
1:C:401:SER:OG	2:C:701:ADP:O2A	2.38	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:216:LEU:HD12	1:E:257:ALA:HB2	2.00	0.42
1:E:554:HIS:CE1	1:F:529:ASP:HB2	2.55	0.42
1:D:60:TYR:CE1	1:D:302:VAL:HG11	2.54	0.42
1:H:202:ILE:HB	1:H:268:LEU:HD13	2.02	0.42
1:C:437:LEU:HD23	1:C:518:ILE:HB	2.00	0.42
1:E:514:LYS:NZ	1:F:236:TYR:O	2.47	0.42
1:A:158:LEU:HG	1:A:325:ILE:HD12	2.02	0.42
1:A:450:THR:HA	1:A:490:GLU:HA	2.01	0.42
1:B:358:ARG:HA	1:B:358:ARG:HD2	1.82	0.42
1:E:181:GLY:O	1:E:182:LEU:HG	2.19	0.42
1:A:243:GLU:OE1	1:A:246:ARG:NH2	2.53	0.42
1:D:375:GLU:HG2	1:D:378:HIS:CD2	2.54	0.42
1:D:437:LEU:HD23	1:D:518:ILE:HB	2.02	0.42
1:E:397:GLY:HA2	2:E:701:ADP:H5'1	2.01	0.42
1:G:237:PHE:CD2	1:H:514:LYS:HD2	2.55	0.42
1:G:566:LEU:HD23	1:G:576:GLN:HG3	2.02	0.42
1:B:50:LYS:HD2	1:B:168:GLU:HB2	2.01	0.41
1:E:221:LEU:HD11	1:F:148:THR:HG21	2.02	0.41
1:G:24:LEU:HD23	1:G:27:LEU:HD12	2.02	0.41
1:B:184:LEU:HD12	1:B:185:VAL:N	2.35	0.41
1:A:493:GLU:OE1	1:B:136:ARG:NH1	2.50	0.41
1:B:223:ARG:HD3	1:B:246:ARG:HG2	2.02	0.41
1:C:262:GLU:HG3	1:D:106:GLU:OE2	2.19	0.41
1:C:261:SER:O	1:C:264:SER:OG	2.37	0.41
1:D:540:MET:HA	1:D:543:VAL:HG22	2.03	0.41
1:G:157:PHE:CD1	1:G:321:VAL:HG22	2.55	0.41
1:H:146:ARG:HH12	1:H:331:ASP:CG	2.28	0.41
1:A:581:ASP:O	1:A:585:ARG:HG3	2.20	0.41
1:A:287:THR:HG21	1:A:301:LEU:HA	2.03	0.41
1:B:22:ARG:NH1	1:D:364:ASP:OD2	2.52	0.41
1:B:313:ARG:HB3	1:B:314:PRO:HD3	2.02	0.41
1:B:346:ALA:HB2	1:B:430:GLN:HE21	1.85	0.41
1:E:429:THR:OG1	1:E:432:SER:OG	2.27	0.41
1:G:130:ARG:HG3	1:G:131:PHE:H	1.86	0.41
1:G:206:ARG:NH1	1:G:265:LEU:HB2	2.36	0.41
1:G:391:ALA:HB3	1:G:565:ILE:HG12	2.03	0.41
1:A:76:ALA:HA	1:A:79:VAL:HG22	2.03	0.41
1:B:61:LYS:HB2	1:B:303:PHE:HB2	2.01	0.41
1:B:579:HIS:ND1	1:B:589:TYR:OH	2.45	0.41
1:C:224:ALA:HB2	1:C:247:TYR:CE1	2.56	0.41
1:E:106:GLU:OE2	1:F:262:GLU:HG3	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:566:LEU:HD23	1:F:576:GLN:HG3	2.03	0.41
1:A:42:ALA:O	1:A:46:VAL:HG23	2.20	0.41
1:H:430:GLN:HB3	1:H:434:ARG:NH1	2.36	0.41
1:A:35:LEU:HD12	1:A:35:LEU:N	2.29	0.41
1:B:22:ARG:NE	1:D:417:ARG:HH21	2.19	0.41
1:B:157:PHE:HD1	1:B:157:PHE:HA	1.75	0.41
1:C:540:MET:HA	1:C:543:VAL:HG22	2.03	0.41
1:D:13:ARG:HA	1:D:13:ARG:HE	1.85	0.41
1:D:35:LEU:HD22	1:D:108:VAL:HG22	2.03	0.41
1:D:41:GLY:O	1:D:45:MET:HG2	2.21	0.41
1:D:446:LEU:HD11	1:D:503:LYS:HG2	2.03	0.41
1:F:54:LEU:O	1:F:57:PRO:HD2	2.21	0.41
1:F:213:MET:HA	1:F:257:ALA:HB1	2.03	0.41
1:F:520:LEU:HD11	1:F:550:ILE:HD12	2.03	0.41
1:G:473:GLY:O	1:G:539:THR:HG21	2.21	0.41
1:H:15:ASP:OD1	1:H:16:GLY:N	2.53	0.41
1:H:113:THR:HG21	1:H:152:ASP:OD1	2.20	0.41
1:H:400:LYS:NZ	2:H:701:ADP:O1B	2.54	0.41
1:H:401:SER:N	2:H:701:ADP:O1A	2.43	0.41
1:D:150:SER:OG	1:D:328:GLY:O	2.37	0.41
1:G:437:LEU:HD23	1:G:518:ILE:HB	2.03	0.41
1:B:317:MET:HE2	1:B:317:MET:HB3	1.91	0.40
1:B:454:ASN:HB3	1:B:510:ARG:HH21	1.86	0.40
1:F:98:ASP:O	1:F:101:ARG:HG2	2.21	0.40
1:C:95:VAL:HG11	1:D:273:ALA:HB2	2.02	0.40
1:C:129:LEU:HD22	1:C:410:PHE:CD2	2.57	0.40
1:C:354:LEU:HD12	1:C:423:GLN:OE1	2.21	0.40
1:C:596:GLN:HG2	1:D:530:THR:HG21	2.03	0.40
1:D:101:ARG:HG3	1:D:160:PHE:HE1	1.86	0.40
1:D:581:ASP:O	1:D:585:ARG:HG3	2.21	0.40
1:E:287:THR:HG21	1:E:301:LEU:HA	2.03	0.40
1:A:216:LEU:HD12	1:A:257:ALA:HB2	2.01	0.40
1:A:458:GLY:HA2	1:B:237:PHE:CD1	2.56	0.40
1:B:157:PHE:O	1:B:161:ASN:HB3	2.22	0.40
1:C:138:GLY:O	1:C:142:LYS:HG2	2.21	0.40
1:G:313:ARG:HB3	1:G:314:PRO:HD3	2.02	0.40
1:G:534:GLN:HG3	1:H:599:GLU:OE2	2.21	0.40
1:A:35:LEU:H	1:A:35:LEU:CD1	2.29	0.40
1:B:217:ASP:HB2	1:B:254:TYR:CE1	2.57	0.40
1:C:457:TYR:OH	1:D:234:VAL:HG22	2.22	0.40
1:E:223:ARG:HD3	1:E:246:ARG:HG2	2.03	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:120:VAL:O	1:G:124:LEU:HD23	2.22	0.40
1:H:42:ALA:O	1:H:46:VAL:HG23	2.21	0.40
1:A:437:LEU:O	1:A:514:LYS:NZ	2.48	0.40
1:E:332:MET:HE3	1:E:332:MET:HB2	1.95	0.40
1:G:229:LEU:HD11	1:H:137:THR:HG23	2.03	0.40

There are no symmetry-related clashes.

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	584/614 (95%)	572 (98%)	11 (2%)	1 (0%)	43	72
1	B	584/614 (95%)	565 (97%)	19 (3%)	0	100	100
1	C	588/614 (96%)	573 (97%)	14 (2%)	1 (0%)	43	72
1	D	584/614 (95%)	572 (98%)	11 (2%)	1 (0%)	43	72
1	E	587/614 (96%)	576 (98%)	11 (2%)	0	100	100
1	F	587/614 (96%)	574 (98%)	12 (2%)	1 (0%)	43	72
1	G	588/614 (96%)	581 (99%)	7 (1%)	0	100	100
1	H	588/614 (96%)	581 (99%)	7 (1%)	0	100	100
All	All	4690/4912 (96%)	4594 (98%)	92 (2%)	4 (0%)	48	78

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	VAL
1	C	182	LEU
1	D	182	LEU
1	F	182	LEU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	469/492 (95%)	467 (100%)	2 (0%)	84	81
1	B	469/492 (95%)	467 (100%)	2 (0%)	84	81
1	C	472/492 (96%)	469 (99%)	3 (1%)	78	79
1	D	469/492 (95%)	467 (100%)	2 (0%)	84	81
1	E	471/492 (96%)	470 (100%)	1 (0%)	87	86
1	F	471/492 (96%)	469 (100%)	2 (0%)	84	81
1	G	472/492 (96%)	470 (100%)	2 (0%)	84	81
1	H	472/492 (96%)	471 (100%)	1 (0%)	87	86
All	All	3765/3936 (96%)	3750 (100%)	15 (0%)	84	81

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	VAL
1	A	153	THR
1	B	108	VAL
1	B	154	MET
1	C	154	MET
1	C	335	MET
1	C	526	SER
1	D	26	TYR
1	D	125	HIS
1	E	26	TYR
1	F	114	ARG
1	F	158	LEU
1	G	125	HIS
1	G	307	TYR
1	H	307	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	208	HIS
1	A	305	ASN
1	A	546	HIS
1	B	18	GLN
1	B	208	HIS
1	B	305	ASN
1	B	327	GLN
1	B	534	GLN
1	B	546	HIS
1	B	596	GLN
1	C	32	ASN
1	C	208	HIS
1	C	327	GLN
1	C	357	ASN
1	C	534	GLN
1	D	14	HIS
1	D	32	ASN
1	D	305	ASN
1	D	327	GLN
1	D	378	HIS
1	D	546	HIS
1	E	208	HIS
1	E	327	GLN
1	E	534	GLN
1	E	546	HIS
1	F	115	HIS
1	F	219	GLN
1	F	310	GLN
1	F	504	GLN
1	F	534	GLN
1	F	546	HIS
1	G	430	GLN
1	G	546	HIS
1	H	365	ASN
1	H	546	HIS
1	H	554	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 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$
2	ADP	F	701	3	28,29,29	1.39	5 (17%)	43,45,45	1.61	9 (20%)
2	ADP	H	701	3	28,29,29	1.39	4 (14%)	43,45,45	1.64	7 (16%)
2	ADP	G	701	3	28,29,29	1.41	5 (17%)	43,45,45	1.62	9 (20%)
2	ADP	D	701	3	28,29,29	1.39	4 (14%)	43,45,45	1.65	8 (18%)
2	ADP	B	701	3	28,29,29	1.41	4 (14%)	43,45,45	1.85	9 (20%)
2	ADP	E	701	3	28,29,29	1.42	4 (14%)	43,45,45	1.84	10 (23%)
2	ADP	A	701	3	28,29,29	1.44	4 (14%)	43,45,45	1.83	9 (20%)
2	ADP	C	701	3	28,29,29	1.42	3 (10%)	43,45,45	1.64	7 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	F	701	3	-	6/16/32/32	0/3/3/3
2	ADP	H	701	3	-	1/16/32/32	0/3/3/3
2	ADP	G	701	3	-	4/16/32/32	0/3/3/3
2	ADP	D	701	3	-	3/16/32/32	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	B	701	3	-	1/16/32/32	0/3/3/3
2	ADP	E	701	3	-	3/16/32/32	0/3/3/3
2	ADP	A	701	3	-	2/16/32/32	0/3/3/3
2	ADP	C	701	3	-	5/16/32/32	0/3/3/3

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	701	ADP	C5-C4	4.83	1.47	1.39
2	A	701	ADP	C5-C4	4.81	1.47	1.39
2	B	701	ADP	C5-C4	4.80	1.47	1.39
2	H	701	ADP	C5-C4	4.74	1.47	1.39
2	E	701	ADP	C5-C4	4.72	1.47	1.39
2	D	701	ADP	C5-C4	4.68	1.47	1.39
2	G	701	ADP	C5-C4	4.63	1.47	1.39
2	F	701	ADP	C5-C4	4.53	1.47	1.39
2	A	701	ADP	C5-C6	2.75	1.48	1.41
2	E	701	ADP	C5-C6	2.74	1.48	1.41
2	B	701	ADP	C5-C6	2.73	1.48	1.41
2	G	701	ADP	C5-C6	2.61	1.48	1.41
2	C	701	ADP	C5-C6	2.58	1.48	1.41
2	H	701	ADP	C5-C6	2.56	1.48	1.41
2	F	701	ADP	C5-C6	2.55	1.48	1.41
2	D	701	ADP	C5-C6	2.52	1.48	1.41
2	E	701	ADP	C8-N7	2.38	1.36	1.31
2	C	701	ADP	C5-N7	-2.36	1.34	1.39
2	H	701	ADP	C5-N7	-2.36	1.34	1.39
2	A	701	ADP	C8-N7	2.36	1.36	1.31
2	B	701	ADP	C8-N7	2.36	1.36	1.31
2	D	701	ADP	C5-N7	-2.30	1.34	1.39
2	G	701	ADP	C8-N7	2.27	1.36	1.31
2	A	701	ADP	C5-N7	-2.27	1.34	1.39
2	E	701	ADP	C5-N7	-2.23	1.35	1.39
2	B	701	ADP	C5-N7	-2.21	1.35	1.39
2	F	701	ADP	C8-N7	2.20	1.35	1.31
2	G	701	ADP	C5-N7	-2.17	1.35	1.39
2	D	701	ADP	C8-N7	2.14	1.35	1.31
2	F	701	ADP	C5-N7	-2.13	1.35	1.39
2	F	701	ADP	C4-N9	-2.11	1.33	1.37
2	H	701	ADP	C8-N7	2.05	1.35	1.31
2	G	701	ADP	C4-N9	-2.04	1.33	1.37

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	ADP	C5-C4-N3	-5.82	118.70	126.72
2	B	701	ADP	C5-C4-N3	-5.80	118.73	126.72
2	E	701	ADP	C5-C4-N3	-5.74	118.82	126.72
2	C	701	ADP	C5-C4-N3	-5.52	119.11	126.72
2	H	701	ADP	C5-C4-N3	-5.32	119.39	126.72
2	D	701	ADP	C5-C4-N3	-5.21	119.54	126.72
2	G	701	ADP	C5-C4-N3	-4.83	120.06	126.72
2	F	701	ADP	C5-C4-N3	-4.75	120.18	126.72
2	A	701	ADP	N3-C4-N9	4.63	135.04	127.17
2	B	701	ADP	N3-C4-N9	4.63	135.04	127.17
2	E	701	ADP	N3-C4-N9	4.56	134.92	127.17
2	C	701	ADP	N3-C4-N9	4.50	134.82	127.17
2	H	701	ADP	N3-C4-N9	4.43	134.70	127.17
2	D	701	ADP	N3-C4-N9	4.37	134.60	127.17
2	F	701	ADP	N3-C4-N9	3.97	133.91	127.17
2	G	701	ADP	N3-C4-N9	3.95	133.89	127.17
2	E	701	ADP	C2-N3-C4	3.69	120.85	111.83
2	B	701	ADP	C2-N3-C4	3.69	120.83	111.83
2	A	701	ADP	C2-N3-C4	3.67	120.78	111.83
2	E	701	ADP	C4-C5-N7	-3.48	106.60	110.58
2	B	701	ADP	C4-C5-N7	-3.44	106.65	110.58
2	A	701	ADP	C4-C5-N7	-3.40	106.70	110.58
2	G	701	ADP	C4-C5-N7	-3.33	106.78	110.58
2	C	701	ADP	C2-N3-C4	3.25	119.77	111.83
2	E	701	ADP	N3-C2-N1	-3.25	123.66	128.58
2	B	701	ADP	N3-C2-N1	-3.21	123.72	128.58
2	F	701	ADP	C4-C5-N7	-3.20	106.93	110.58
2	F	701	ADP	C4-N9-C8	3.17	109.07	105.74
2	C	701	ADP	C4-C5-N7	-3.16	106.97	110.58
2	A	701	ADP	N3-C2-N1	-3.15	123.81	128.58
2	D	701	ADP	C2-N3-C4	3.15	119.53	111.83
2	H	701	ADP	C2-N3-C4	3.15	119.53	111.83
2	D	701	ADP	C4-C5-N7	-3.09	107.05	110.58
2	H	701	ADP	C4-C5-N7	-3.09	107.05	110.58
2	G	701	ADP	C2-N3-C4	3.08	119.35	111.83
2	F	701	ADP	C2-N3-C4	3.07	119.33	111.83
2	G	701	ADP	C4-N9-C8	3.02	108.91	105.74
2	D	701	ADP	C4-N9-C8	2.76	108.63	105.74
2	E	701	ADP	C4-N9-C8	2.73	108.60	105.74
2	F	701	ADP	N3-C2-N1	-2.72	124.47	128.58
2	B	701	ADP	C4-N9-C8	2.68	108.55	105.74
2	G	701	ADP	N3-C2-N1	-2.64	124.58	128.58

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	ADP	C4-N9-C8	2.62	108.49	105.74
2	E	701	ADP	C5-N7-C8	2.57	107.49	103.45
2	D	701	ADP	N3-C2-N1	-2.53	124.75	128.58
2	B	701	ADP	C5-N7-C8	2.53	107.42	103.45
2	H	701	ADP	C4-N9-C8	2.52	108.38	105.74
2	A	701	ADP	C5-N7-C8	2.49	107.36	103.45
2	C	701	ADP	N3-C2-N1	-2.48	124.83	128.58
2	H	701	ADP	N3-C2-N1	-2.44	124.88	128.58
2	G	701	ADP	C5-N7-C8	2.43	107.27	103.45
2	F	701	ADP	C5-N7-C8	2.38	107.19	103.45
2	D	701	ADP	C5-N7-C8	2.37	107.18	103.45
2	B	701	ADP	C3'-C2'-C1'	2.37	105.94	101.46
2	E	701	ADP	C3'-C2'-C1'	2.35	105.91	101.46
2	A	701	ADP	C3'-C2'-C1'	2.33	105.88	101.46
2	H	701	ADP	C5-N7-C8	2.29	107.05	103.45
2	C	701	ADP	C5-N7-C8	2.28	107.03	103.45
2	C	701	ADP	C4-N9-C8	2.22	108.07	105.74
2	E	701	ADP	C6-C5-N7	2.16	136.26	132.09
2	G	701	ADP	C6-C5-N7	2.15	136.23	132.09
2	F	701	ADP	N9-C8-N7	-2.14	110.91	113.94
2	F	701	ADP	C6-C5-N7	2.10	136.14	132.09
2	B	701	ADP	C6-C5-N7	2.09	136.12	132.09
2	G	701	ADP	N9-C8-N7	-2.08	110.99	113.94
2	E	701	ADP	N9-C8-N7	-2.04	111.04	113.94
2	A	701	ADP	C6-C5-N7	2.03	136.00	132.09
2	D	701	ADP	C2'-C1'-N9	-2.01	108.32	113.30

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	701	ADP	PA-O3A-PB-O2B
2	C	701	ADP	PA-O3A-PB-O3B
2	C	701	ADP	C5'-O5'-PA-O1A
2	D	701	ADP	C5'-O5'-PA-O1A
2	D	701	ADP	C5'-O5'-PA-O2A
2	D	701	ADP	C5'-O5'-PA-O3A
2	E	701	ADP	C5'-O5'-PA-O1A
2	E	701	ADP	C5'-O5'-PA-O2A
2	E	701	ADP	C5'-O5'-PA-O3A
2	F	701	ADP	C5'-O5'-PA-O1A
2	F	701	ADP	C5'-O5'-PA-O2A

Continued on next page...

Continued from previous page...

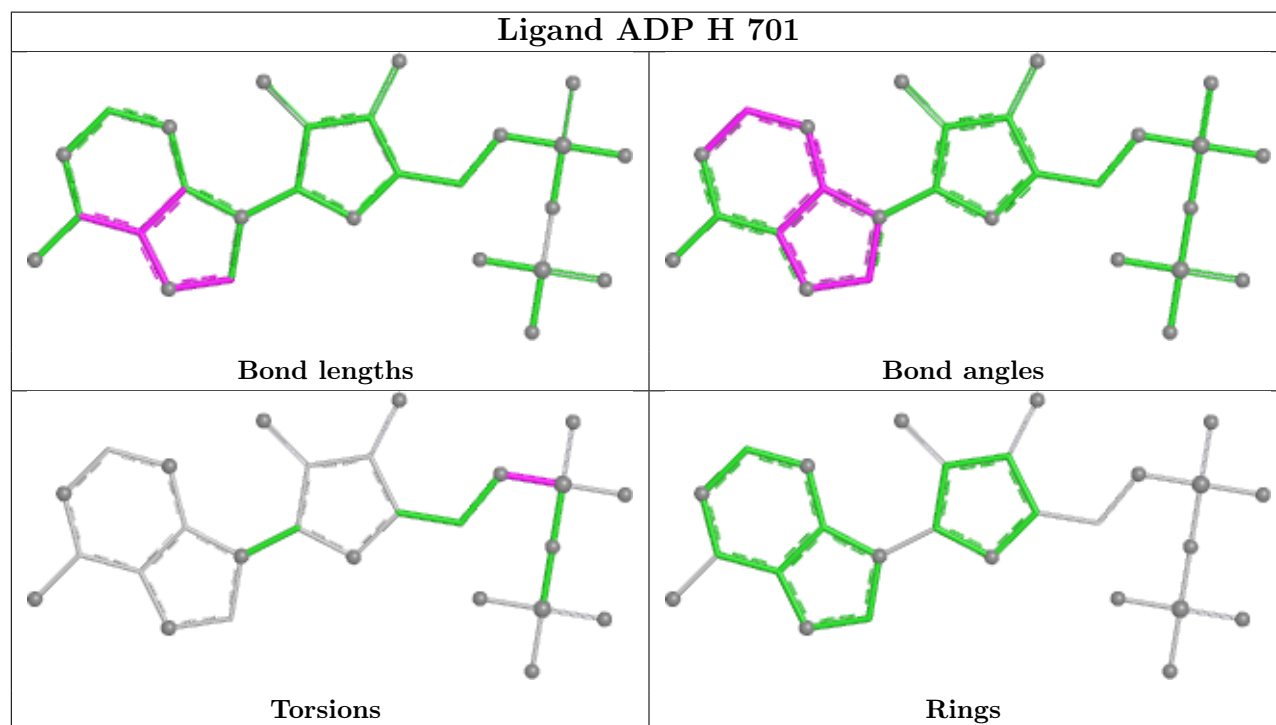
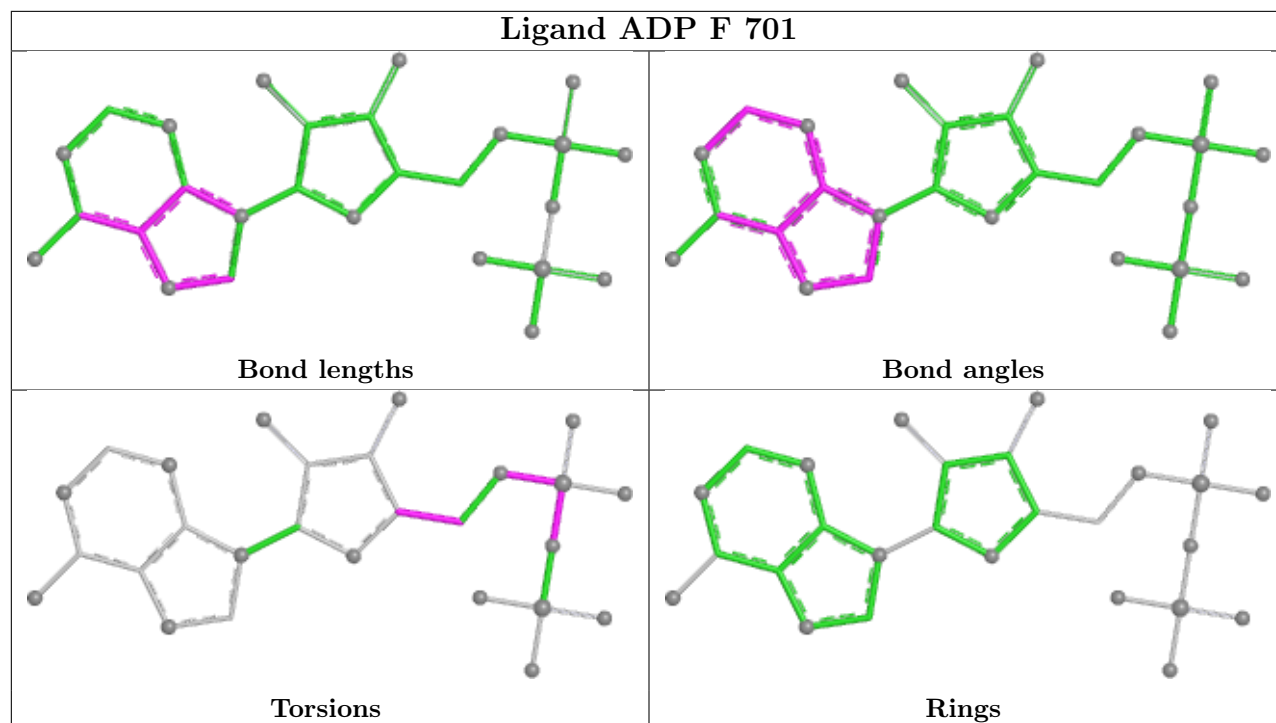
Mol	Chain	Res	Type	Atoms
2	F	701	ADP	C5'-O5'-PA-O3A
2	G	701	ADP	C5'-O5'-PA-O3A
2	F	701	ADP	C3'-C4'-C5'-O5'
2	F	701	ADP	O4'-C4'-C5'-O5'
2	G	701	ADP	O4'-C4'-C5'-O5'
2	C	701	ADP	O4'-C4'-C5'-O5'
2	G	701	ADP	C3'-C4'-C5'-O5'
2	A	701	ADP	C5'-O5'-PA-O1A
2	G	701	ADP	C5'-O5'-PA-O1A
2	H	701	ADP	C5'-O5'-PA-O1A
2	C	701	ADP	PA-O3A-PB-O1B
2	B	701	ADP	O4'-C4'-C5'-O5'
2	A	701	ADP	O4'-C4'-C5'-O5'
2	F	701	ADP	PB-O3A-PA-O2A

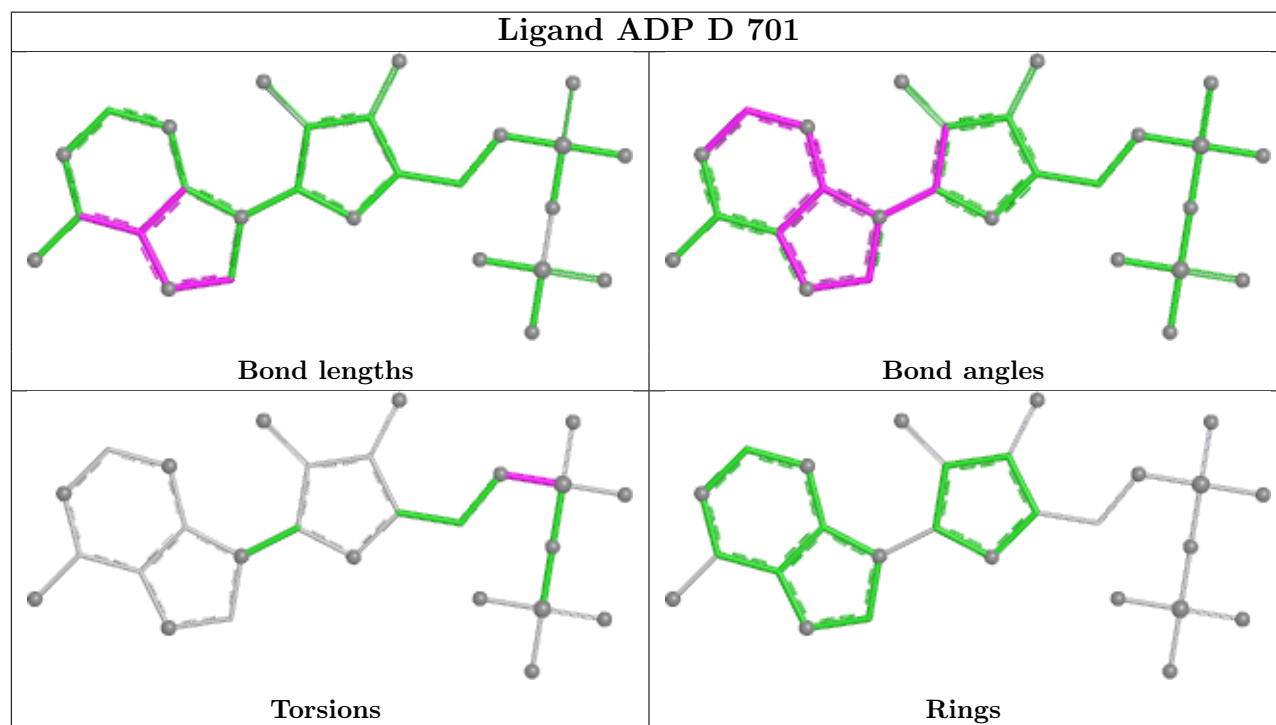
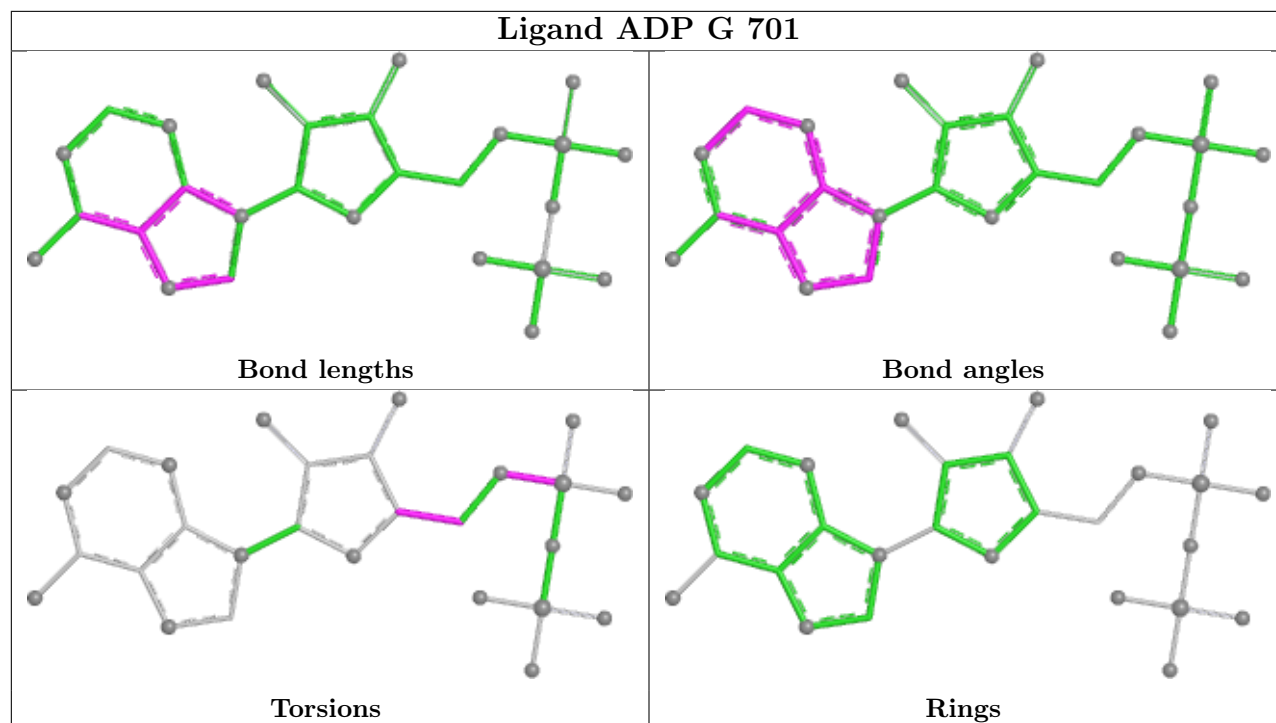
There are no ring outliers.

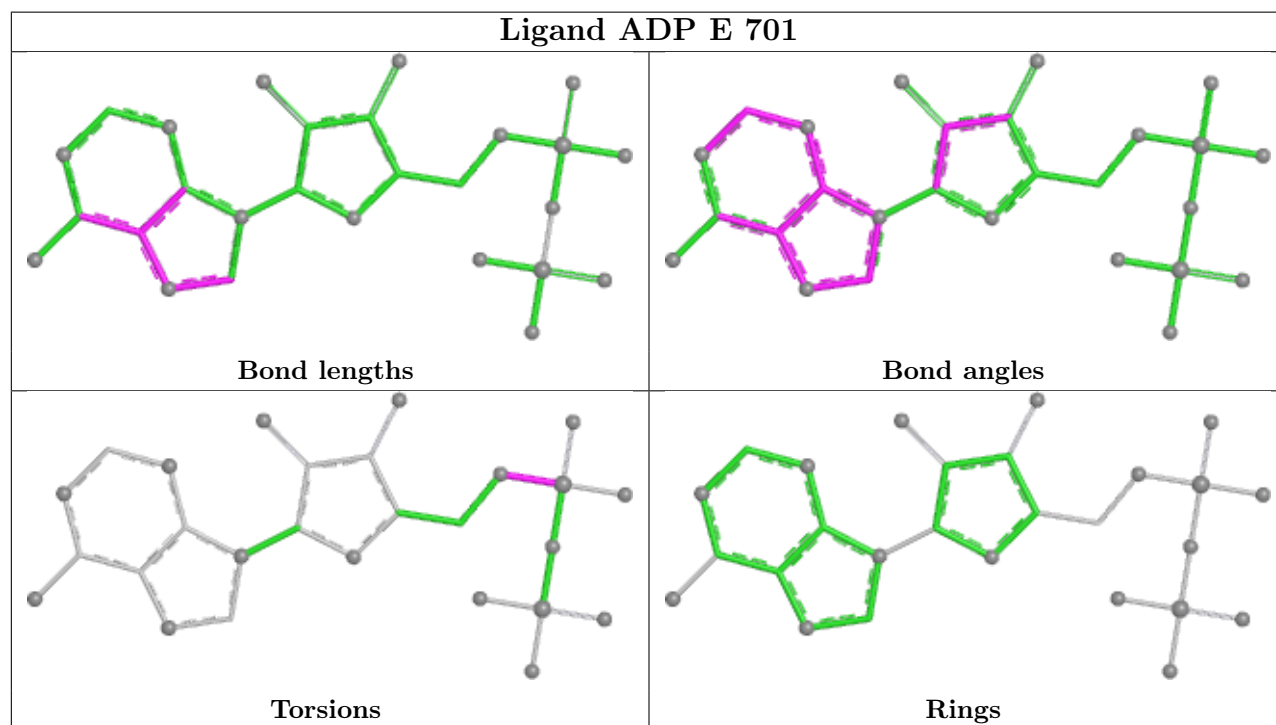
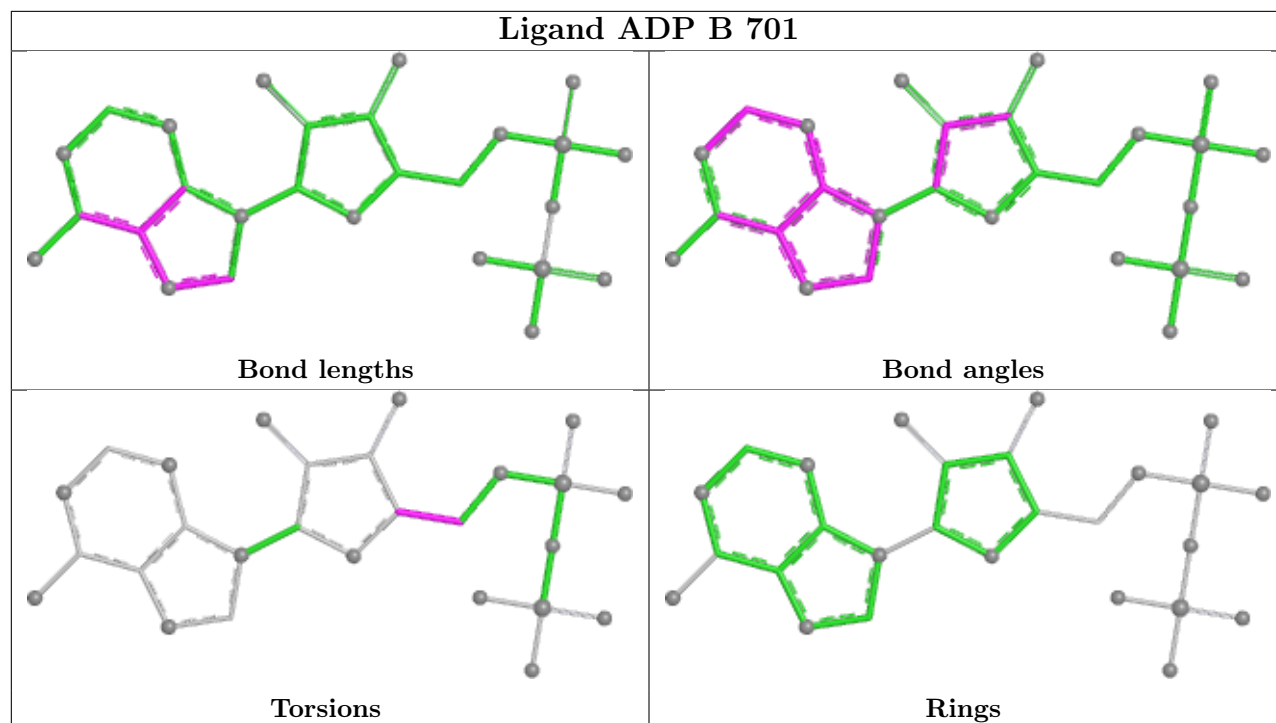
7 monomers are involved in 13 short contacts:

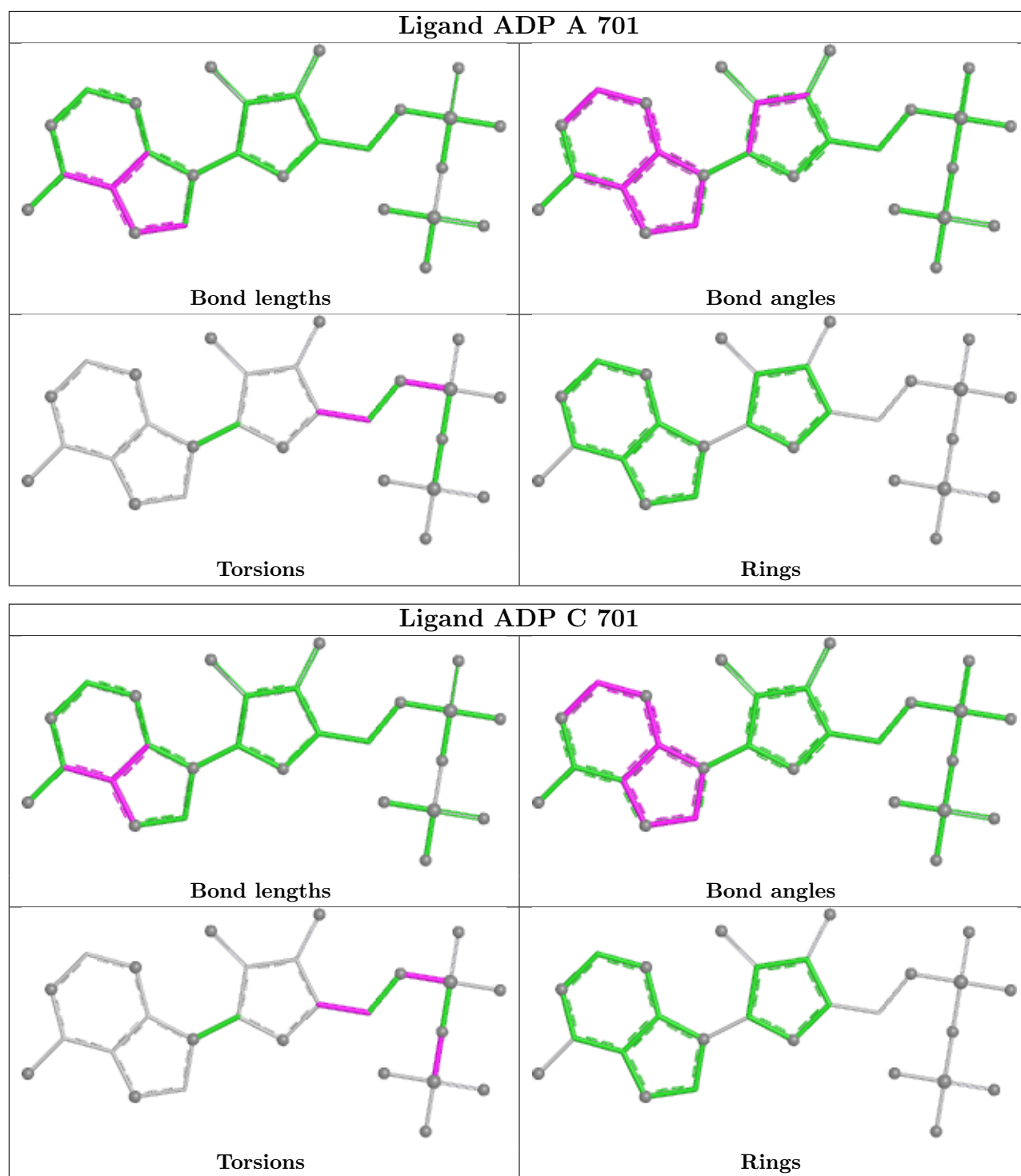
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	701	ADP	1	0
2	H	701	ADP	2	0
2	G	701	ADP	1	0
2	B	701	ADP	2	0
2	E	701	ADP	3	0
2	A	701	ADP	2	0
2	C	701	ADP	2	0

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 ⓘ

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	586/614 (95%)	0.01	19 (3%)	50	31	82, 131, 179, 338	0
1	B	586/614 (95%)	0.02	21 (3%)	46	29	87, 138, 177, 227	0
1	C	590/614 (96%)	-0.03	22 (3%)	45	28	95, 136, 181, 240	0
1	D	586/614 (95%)	0.03	27 (4%)	37	24	98, 142, 186, 400	0
1	E	589/614 (95%)	0.07	24 (4%)	41	26	138, 210, 250, 271	0
1	F	589/614 (95%)	0.25	35 (5%)	28	19	136, 211, 256, 278	0
1	G	590/614 (96%)	0.16	19 (3%)	50	31	158, 224, 253, 262	0
1	H	590/614 (96%)	0.25	34 (5%)	29	20	179, 217, 251, 270	0
All	All	4706/4912 (95%)	0.09	201 (4%)	40	25	82, 176, 245, 400	0

All (201) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	67	MET	6.2
1	F	563	ASP	6.0
1	D	139	GLU	5.7
1	D	70	GLY	5.4
1	D	132	HIS	5.1
1	F	231	TYR	5.0
1	B	356	VAL	4.8
1	B	74	GLN	4.8
1	D	135	ARG	4.8
1	D	138	GLY	4.7
1	B	300	ASP	4.7
1	B	353	ALA	4.7
1	A	163	ALA	4.6
1	H	445	VAL	4.6
1	H	348	VAL	4.4
1	E	535	ASP	4.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	597	ALA	4.3
1	A	27	LEU	4.3
1	D	137	THR	4.2
1	E	125	HIS	4.2
1	F	386	ALA	4.1
1	G	488	ASP	3.9
1	E	478	ASP	3.7
1	D	371	ASP	3.7
1	G	520	LEU	3.7
1	F	562	SER	3.6
1	F	238	GLY	3.6
1	F	389	ARG	3.6
1	D	375	GLU	3.6
1	C	67	MET	3.6
1	H	177	TRP	3.6
1	E	574	ALA	3.5
1	F	416	GLY	3.5
1	H	589	TYR	3.5
1	F	388	SER	3.4
1	H	349	PRO	3.4
1	G	228	LEU	3.4
1	E	569	ASP	3.4
1	D	69	LEU	3.3
1	H	490	GLU	3.3
1	H	514	LYS	3.3
1	E	435	ALA	3.2
1	H	136	ARG	3.2
1	E	396	SER	3.2
1	D	498	LEU	3.2
1	F	58	PHE	3.2
1	A	139	GLU	3.2
1	C	598	ALA	3.2
1	E	76	ALA	3.2
1	B	494	ARG	3.1
1	C	294	GLY	3.1
1	E	86	ALA	3.1
1	E	131	PHE	3.1
1	G	110	GLN	3.1
1	F	387	GLY	3.1
1	H	346	ALA	3.1
1	E	588	LEU	3.1
1	F	383	GLU	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	F	385	ALA	3.0
1	C	494	ARG	3.0
1	H	174	VAL	3.0
1	C	214	ASN	3.0
1	B	354	LEU	3.0
1	G	298	VAL	3.0
1	H	599	GLU	2.9
1	D	60	TYR	2.9
1	C	292	SER	2.9
1	B	355	VAL	2.9
1	A	35	LEU	2.9
1	H	590	ALA	2.9
1	F	35	LEU	2.9
1	H	178	LEU	2.8
1	G	306	THR	2.8
1	A	69	LEU	2.8
1	F	124	LEU	2.8
1	F	301	LEU	2.8
1	C	293	GLN	2.8
1	H	386	ALA	2.8
1	F	110	GLN	2.8
1	F	347	ASP	2.8
1	C	498	LEU	2.8
1	D	373	ASP	2.8
1	C	527	CYS	2.8
1	H	125	HIS	2.7
1	H	60	TYR	2.7
1	E	255	ALA	2.7
1	A	34	VAL	2.7
1	A	95	VAL	2.7
1	E	124	LEU	2.7
1	H	520	LEU	2.7
1	H	588	LEU	2.7
1	G	576	GLN	2.7
1	B	293	GLN	2.6
1	B	597	ALA	2.6
1	H	300	ASP	2.6
1	A	161	ASN	2.6
1	E	531	ARG	2.6
1	C	442	GLN	2.6
1	H	72	GLY	2.6
1	E	418	ILE	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	F	149	LYS	2.6
1	A	371	ASP	2.6
1	B	145	GLU	2.6
1	D	358	ARG	2.6
1	D	73	ALA	2.6
1	F	377	LEU	2.6
1	C	291	TRP	2.5
1	H	446	LEU	2.5
1	C	334	GLU	2.5
1	H	474	ALA	2.5
1	A	318	LEU	2.5
1	A	316	ASP	2.5
1	F	59	ALA	2.5
1	C	210	ARG	2.5
1	A	496	LEU	2.5
1	D	378	HIS	2.5
1	D	149	LYS	2.5
1	B	313	ARG	2.5
1	C	443	ASP	2.5
1	C	161	ASN	2.4
1	B	374	ARG	2.4
1	B	574	ALA	2.4
1	E	434	ARG	2.4
1	B	525	THR	2.4
1	H	600	SER	2.4
1	D	356	VAL	2.4
1	H	30	ALA	2.4
1	G	276	VAL	2.4
1	F	343	ILE	2.4
1	F	60	TYR	2.4
1	E	476	ILE	2.4
1	F	236	TYR	2.4
1	E	130	ARG	2.4
1	F	284	MET	2.4
1	B	359	PRO	2.4
1	G	176	PHE	2.4
1	B	420	ILE	2.4
1	D	141	THR	2.4
1	D	65	ASP	2.4
1	A	67	MET	2.4
1	G	178	LEU	2.3
1	C	60	TYR	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	G	581	ASP	2.3
1	C	211	GLU	2.3
1	F	51	ALA	2.3
1	F	565	ILE	2.3
1	F	55	ALA	2.3
1	B	269	ASN	2.3
1	F	211	GLU	2.3
1	G	304	VAL	2.3
1	G	530	THR	2.3
1	H	179	ASN	2.3
1	F	353	ALA	2.3
1	G	312	PHE	2.3
1	C	145	GLU	2.3
1	D	71	GLY	2.3
1	D	196	VAL	2.2
1	E	395	PRO	2.2
1	G	577	GLY	2.2
1	E	47	LEU	2.2
1	H	586	ASP	2.2
1	C	318	LEU	2.2
1	D	21	LYS	2.2
1	C	108	VAL	2.2
1	D	357	ASN	2.2
1	E	66	ALA	2.2
1	E	73	ALA	2.2
1	G	445	VAL	2.2
1	H	140	VAL	2.2
1	H	144	ILE	2.2
1	H	439	ILE	2.2
1	H	592	MET	2.2
1	H	522	ASP	2.1
1	B	302	VAL	2.1
1	C	493	GLU	2.1
1	A	24	LEU	2.1
1	G	224	ALA	2.1
1	D	126	LYS	2.1
1	E	90	GLY	2.1
1	G	111	ASP	2.1
1	F	111	ASP	2.1
1	B	492	GLY	2.1
1	H	337	ARG	2.1
1	B	295	LYS	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	F	31	ASP	2.1
1	D	68	THR	2.1
1	F	33	ALA	2.1
1	E	311	LEU	2.1
1	A	71	GLY	2.0
1	A	173	ILE	2.0
1	H	444	SER	2.0
1	F	339	ILE	2.0
1	F	595	ARG	2.0
1	A	165	THR	2.0
1	A	159	LEU	2.0
1	D	570	GLN	2.0
1	B	486	GLY	2.0
1	G	287	THR	2.0
1	F	228	LEU	2.0
1	A	444	SER	2.0
1	H	76	ALA	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.

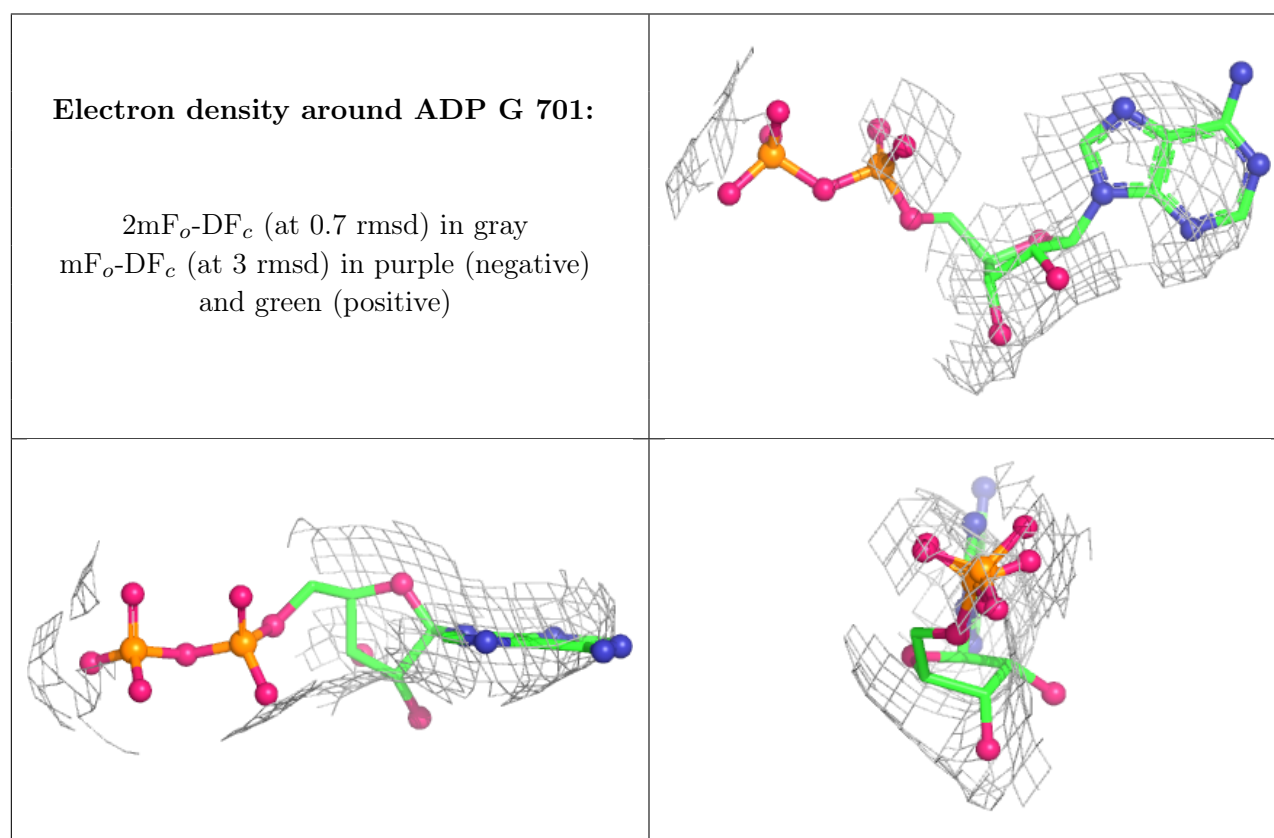
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	D	702	1/1	0.73	0.31	150,150,150,150	0
2	ADP	G	701	27/27	0.82	0.10	241,255,282,283	0
2	ADP	C	701	27/27	0.83	0.12	172,188,203,206	0
2	ADP	E	701	27/27	0.84	0.12	174,183,194,197	0
2	ADP	D	701	27/27	0.85	0.13	154,184,200,205	0
2	ADP	F	701	27/27	0.85	0.08	196,203,210,212	0

Continued on next page...

Continued from previous page...

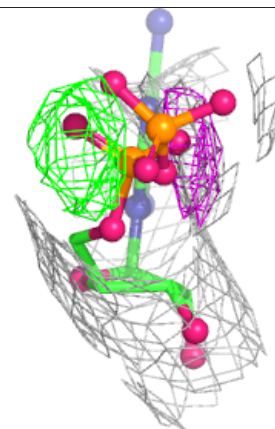
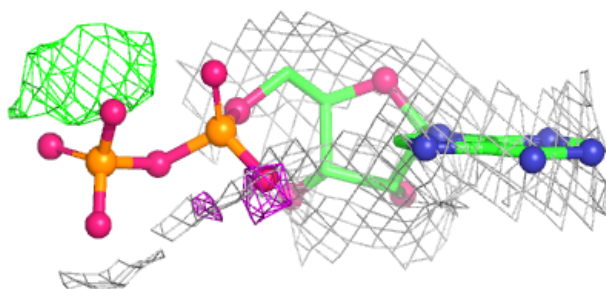
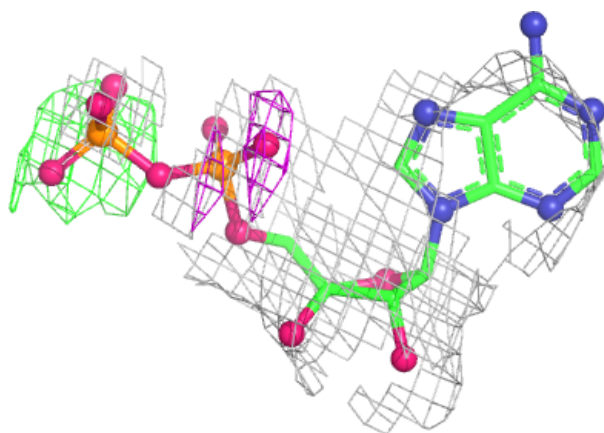
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ADP	H	701	27/27	0.86	0.12	213,222,226,229	0
2	ADP	B	701	27/27	0.87	0.12	159,183,220,226	0
3	MG	G	702	1/1	0.88	0.11	151,151,151,151	0
3	MG	C	702	1/1	0.92	0.13	89,89,89,89	0
2	ADP	A	701	27/27	0.93	0.11	123,144,170,175	0
3	MG	F	702	1/1	0.94	0.18	120,120,120,120	0
3	MG	B	702	1/1	0.94	0.16	92,92,92,92	0
3	MG	E	702	1/1	0.95	0.14	99,99,99,99	0
3	MG	H	702	1/1	0.97	0.17	138,138,138,138	0
3	MG	A	702	1/1	0.98	0.07	74,74,74,74	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.

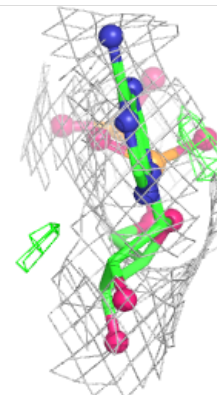
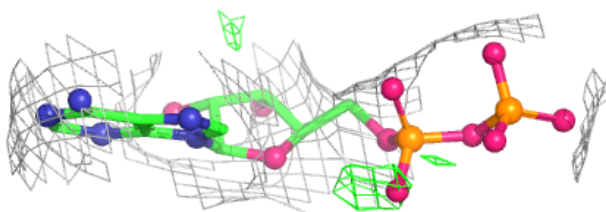
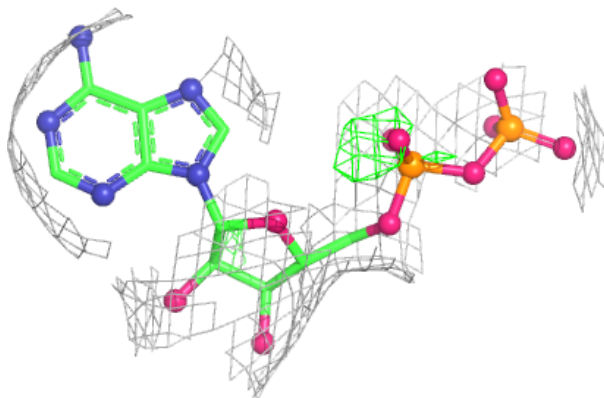


Electron density around ADP C 701:

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

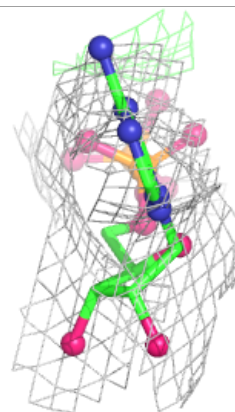
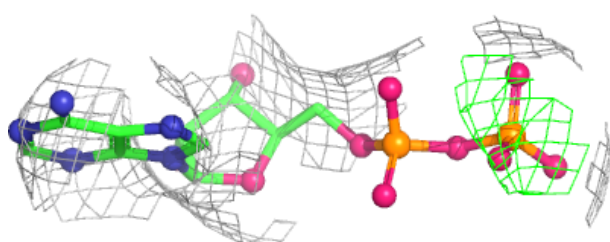
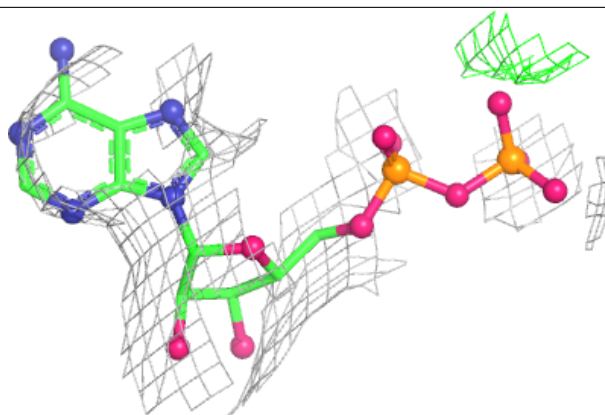
**Electron density around ADP E 701:**

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

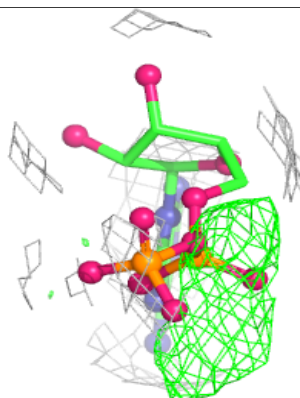
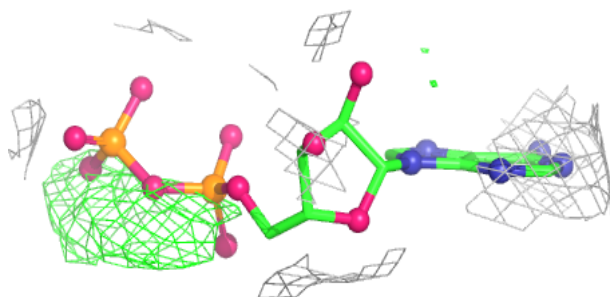
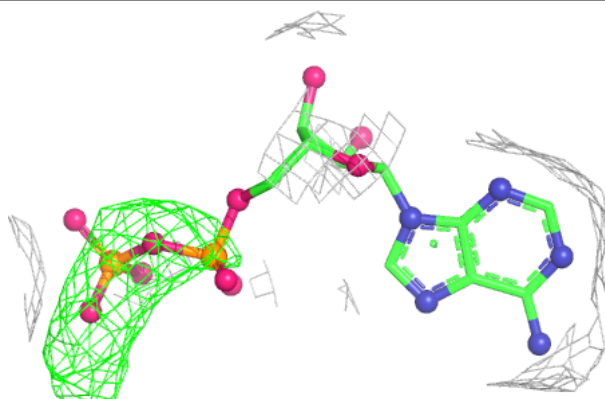


Electron density around ADP D 701:

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

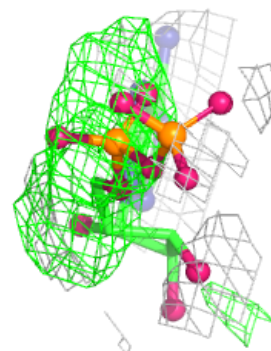
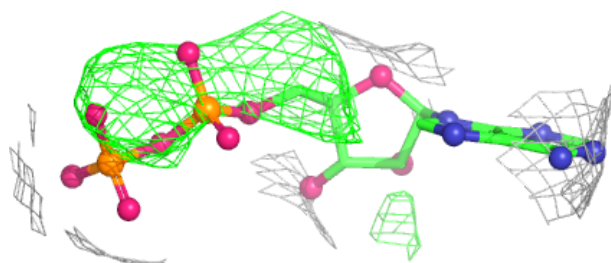
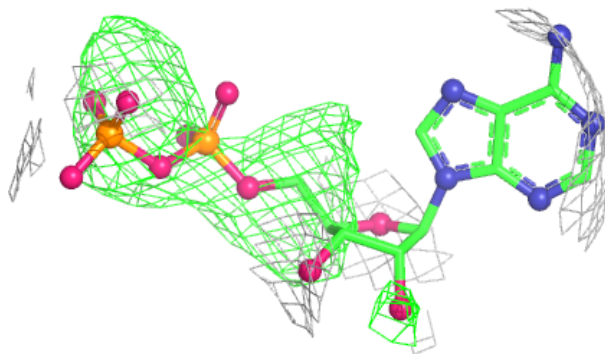
**Electron density around ADP F 701:**

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

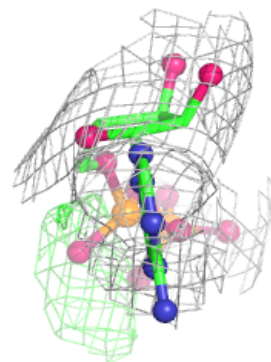
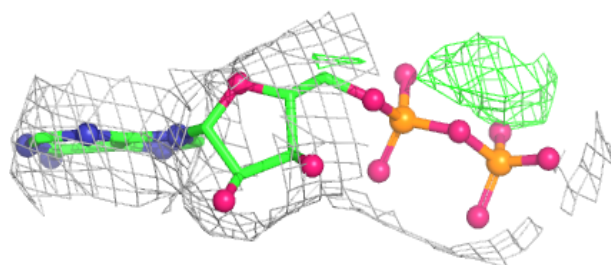
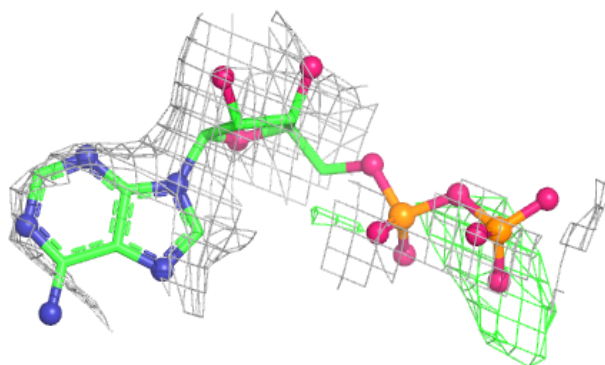


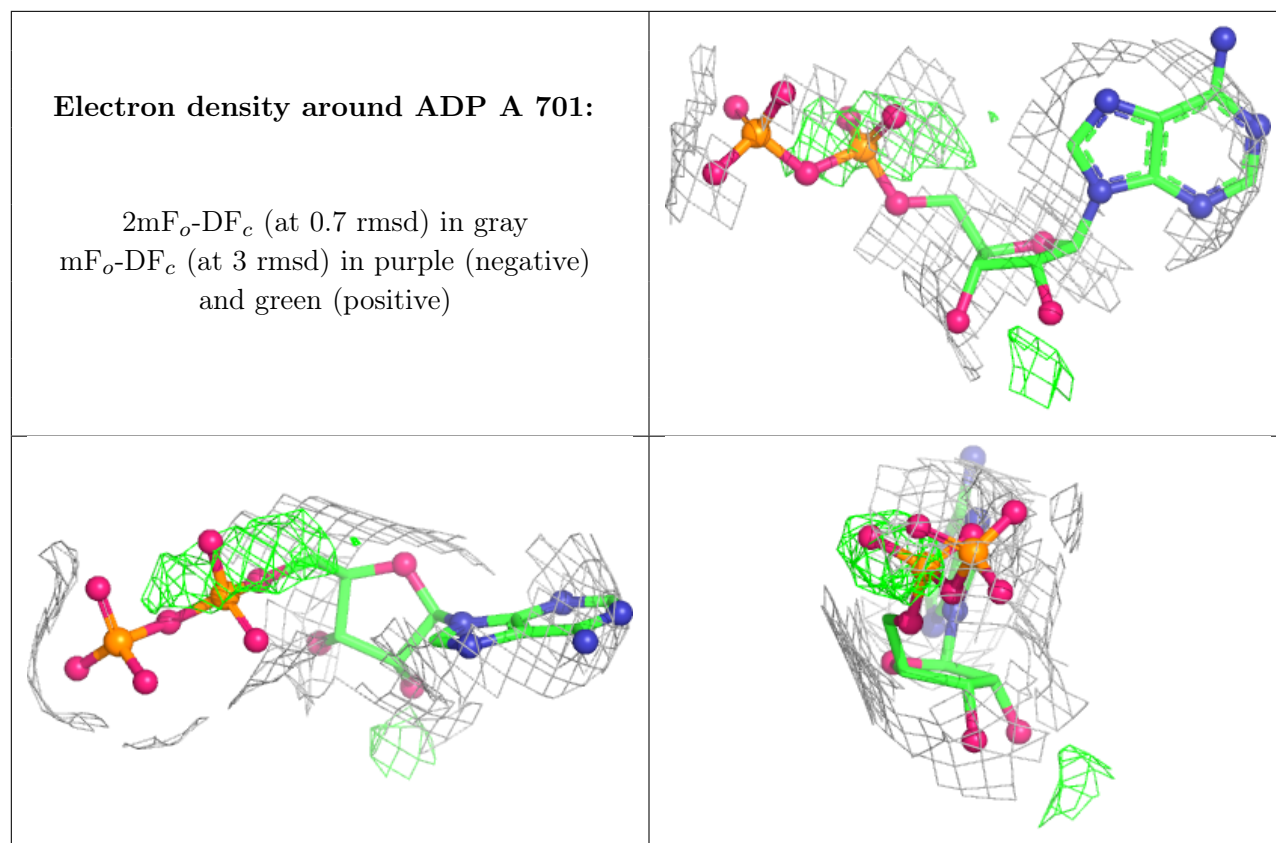
Electron density around ADP H 701:

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

**Electron density around ADP B 701:**

$2mF_o-DF_c$ (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.