



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:25 AM UTC

PDB ID : 4IC7 / pdb_00004ic7
Title : Crystal structure of the ERK5 kinase domain in complex with an MKK5 binding fragment
Authors : Gogl, G.; Remenyi, A.
Deposited on : 2012-12-10
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

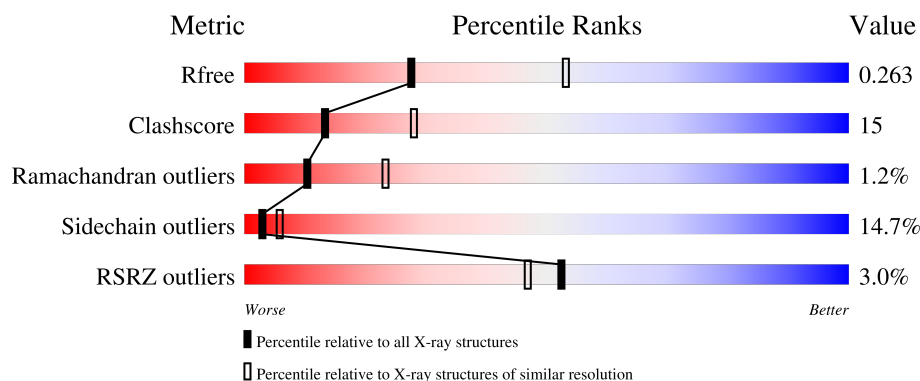
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	442	2% 51% 24% 5% 20%
1	D	442	2% 52% 21% 6% 20%
2	B	126	4% 46% 37% 12% ..
2	E	126	4% 55% 28% 5% • 12%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7394 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 Mitogen-activated protein kinase 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	355	Total	C	N	O	S	0	0	0
			2775	1776	488	498	13			
1	D	353	Total	C	N	O	S	0	0	0
			2675	1713	471	478	13			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q13164
A	0	SER	-	expression tag	UNP Q13164
A	432	SER	-	expression tag	UNP Q13164
A	433	GLY	-	expression tag	UNP Q13164
A	434	ARG	-	expression tag	UNP Q13164
A	435	HIS	-	expression tag	UNP Q13164
A	436	HIS	-	expression tag	UNP Q13164
A	437	HIS	-	expression tag	UNP Q13164
A	438	HIS	-	expression tag	UNP Q13164
A	439	HIS	-	expression tag	UNP Q13164
A	440	HIS	-	expression tag	UNP Q13164
D	-1	GLY	-	expression tag	UNP Q13164
D	0	SER	-	expression tag	UNP Q13164
D	432	SER	-	expression tag	UNP Q13164
D	433	GLY	-	expression tag	UNP Q13164
D	434	ARG	-	expression tag	UNP Q13164
D	435	HIS	-	expression tag	UNP Q13164
D	436	HIS	-	expression tag	UNP Q13164
D	437	HIS	-	expression tag	UNP Q13164
D	438	HIS	-	expression tag	UNP Q13164
D	439	HIS	-	expression tag	UNP Q13164
D	440	HIS	-	expression tag	UNP Q13164

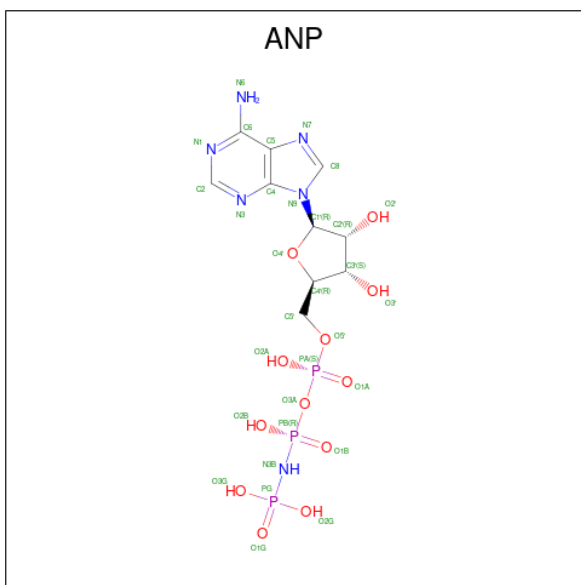
- Molecule 2 is a protein called Dual specificity mitogen-activated protein kinase kinase 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	111	Total	C	N	O	S	0	0	0
			859	544	145	166	4			
2	B	123	Total	C	N	O	S	0	0	0
			940	592	166	178	4			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	14	GLY	-	expression tag	UNP Q13163
E	15	SER	-	expression tag	UNP Q13163
E	131	SER	-	expression tag	UNP Q13163
E	132	GLY	-	expression tag	UNP Q13163
E	133	ARG	-	expression tag	UNP Q13163
E	134	HIS	-	expression tag	UNP Q13163
E	135	HIS	-	expression tag	UNP Q13163
E	136	HIS	-	expression tag	UNP Q13163
E	137	HIS	-	expression tag	UNP Q13163
E	138	HIS	-	expression tag	UNP Q13163
E	139	HIS	-	expression tag	UNP Q13163
B	14	GLY	-	expression tag	UNP Q13163
B	15	SER	-	expression tag	UNP Q13163
B	131	SER	-	expression tag	UNP Q13163
B	132	GLY	-	expression tag	UNP Q13163
B	133	ARG	-	expression tag	UNP Q13163
B	134	HIS	-	expression tag	UNP Q13163
B	135	HIS	-	expression tag	UNP Q13163
B	136	HIS	-	expression tag	UNP Q13163
B	137	HIS	-	expression tag	UNP Q13163
B	138	HIS	-	expression tag	UNP Q13163
B	139	HIS	-	expression tag	UNP Q13163

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 31	C 10	N 6	O 12	P 3	0	0
3	D	1	Total 31	C 10	N 6	O 12	P 3	0	0

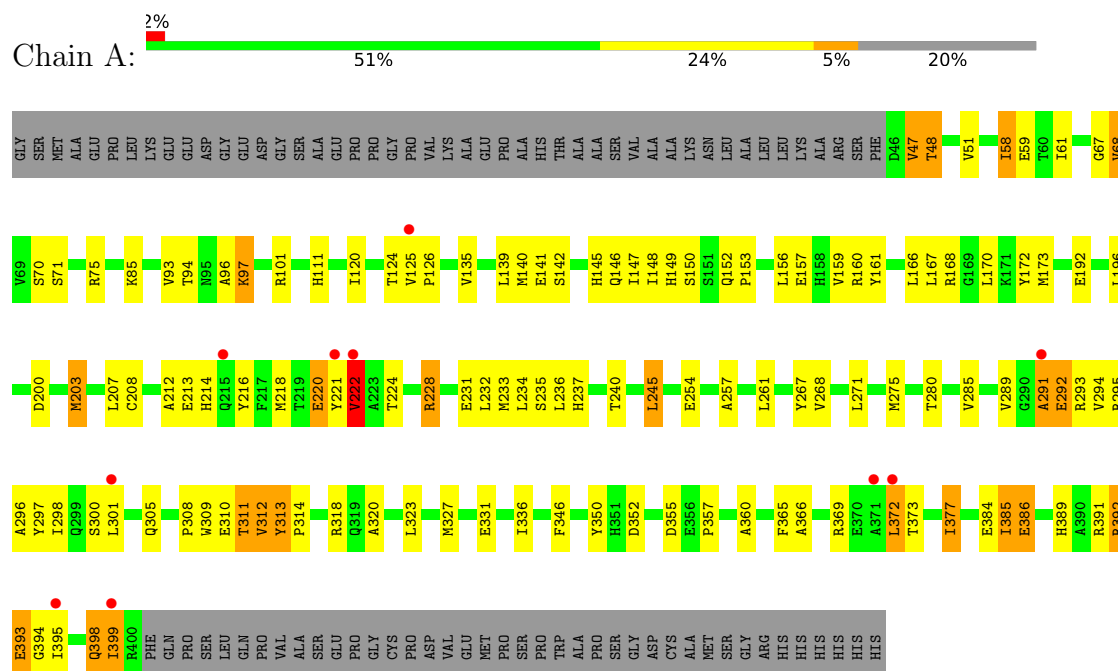
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	36	Total O 36 36	0	0
4	D	32	Total O 32 32	0	0
4	E	10	Total O 10 10	0	0
4	B	5	Total O 5 5	0	0

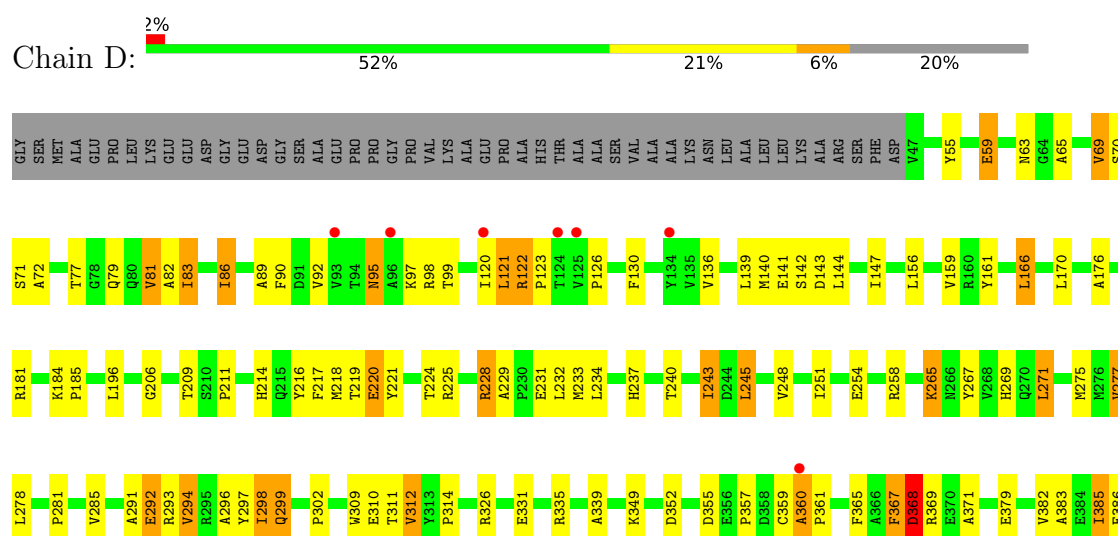
3 Residue-property plots

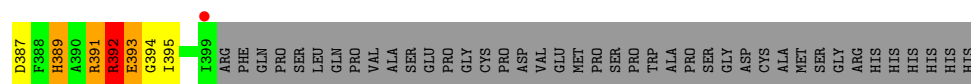
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Mitogen-activated protein kinase 7

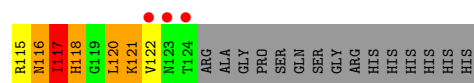


• Molecule 1: Mitogen-activated protein kinase 7

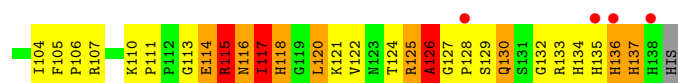
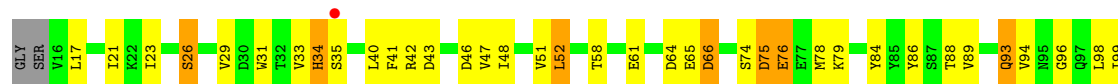




- Molecule 2: Dual specificity mitogen-activated protein kinase kinase 5



- Molecule 2: Dual specificity mitogen-activated protein kinase kinase 5



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	69.37Å 69.37Å 271.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.27 – 2.60 48.27 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.6 (48.27-2.60) 94.5 (48.27-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.211 , 0.258 0.217 , 0.263	Depositor DCC
R_{free} test set	1958 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	56.8	Xtriage
Anisotropy	0.613	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 76.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.104 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7394	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.28% 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: ANP

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	0.57	0/2843	1.09	15/3865 (0.4%)
1	D	0.54	0/2740	1.09	17/3735 (0.5%)
2	B	0.63	0/961	1.18	9/1311 (0.7%)
2	E	0.57	0/876	1.02	4/1194 (0.3%)
All	All	0.57	0/7420	1.10	45/10105 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	D	0	3
2	B	0	5
2	E	0	1
All	All	0	12

There are no bond length outliers.

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	360	ALA	CA-C-N	-12.01	108.01	120.38
1	D	360	ALA	C-N-CA	-12.01	108.01	120.38
1	A	360	ALA	CA-C-N	-11.83	108.19	120.38
1	A	360	ALA	C-N-CA	-11.83	108.19	120.38
2	E	118	HIS	N-CA-C	9.91	125.57	109.72
2	B	118	HIS	N-CA-C	9.68	125.14	109.85
1	D	294	VAL	N-CA-C	-9.03	99.35	111.44
1	A	292	GLU	N-CA-C	7.59	126.97	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	121	LYS	N-CA-C	7.36	120.00	111.02
1	A	291	ALA	N-CA-C	7.10	119.66	108.79
1	D	392	ARG	CA-C-N	7.06	134.41	121.70
1	D	392	ARG	C-N-CA	7.06	134.41	121.70
1	D	86	ILE	CA-C-N	6.79	128.33	119.84
1	D	86	ILE	C-N-CA	6.79	128.33	119.84
1	A	267	TYR	N-CA-C	6.69	118.26	110.97
2	B	125	ARG	N-CA-C	-6.55	100.51	110.14
2	B	110	LYS	CA-C-N	6.49	127.07	120.38
2	B	110	LYS	C-N-CA	6.49	127.07	120.38
1	A	67	GLY	N-CA-C	6.32	118.47	110.96
2	B	117	ILE	N-CA-C	6.30	122.44	109.34
1	A	394	GLY	N-CA-C	-6.29	98.27	113.18
1	A	386	GLU	N-CA-C	-6.28	103.11	111.96
1	A	294	VAL	N-CA-C	-6.22	103.86	112.50
1	D	122	ARG	CA-C-N	6.17	127.56	119.84
1	D	122	ARG	C-N-CA	6.17	127.56	119.84
1	A	214	HIS	N-CA-C	-6.17	100.17	109.79
2	E	66	ASP	N-CA-C	-6.02	105.94	113.28
1	D	297	TYR	N-CA-C	-5.94	98.14	110.80
2	B	75	ASP	N-CA-C	-5.92	100.54	109.60
1	D	292	GLU	N-CA-C	5.89	123.35	110.80
1	A	222	VAL	N-CA-C	-5.77	97.35	109.34
1	D	267	TYR	N-CA-C	5.64	117.88	111.11
2	E	117	ILE	N-CA-C	5.50	120.79	109.34
2	B	126	ALA	N-CA-C	5.41	122.32	110.80
2	B	52	LEU	CA-C-N	5.39	124.90	119.19
2	B	52	LEU	C-N-CA	5.39	124.90	119.19
1	A	207	LEU	CB-CA-C	5.36	121.09	110.42
1	D	229	ALA	CA-C-N	5.17	125.21	119.32
1	D	229	ALA	C-N-CA	5.17	125.21	119.32
1	A	298	ILE	N-CA-C	-5.16	105.81	110.82
1	A	313	TYR	CA-C-N	5.08	124.25	118.97
1	A	313	TYR	C-N-CA	5.08	124.25	118.97
1	D	357	PRO	N-CA-C	5.05	119.41	111.38
1	D	302	PRO	CA-C-N	5.04	125.04	119.90
1	D	302	PRO	C-N-CA	5.04	125.04	119.90

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	292	GLU	Peptide
1	A	366	ALA	Peptide
1	A	51	VAL	Peptide
2	B	115	ARG	Sidechain
2	B	117	ILE	Peptide
2	B	126	ALA	Peptide
2	B	130	GLN	Peptide
2	B	136	HIS	Peptide
1	D	292	GLU	Peptide
1	D	367	PHE	Peptide
1	D	392	ARG	Peptide
2	E	117	ILE	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2775	0	2698	77	0
1	D	2675	0	2551	70	0
2	B	940	0	891	53	0
2	E	859	0	828	30	0
3	A	31	0	11	3	0
3	D	31	0	13	8	0
4	A	36	0	0	4	0
4	B	5	0	0	1	0
4	D	32	0	0	2	0
4	E	10	0	0	0	0
All	All	7394	0	6992	212	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 15.

All (212) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:THR:O	1:A:85:LYS:NZ	1.99	0.95
2:E:112:PRO:HA	2:E:114:GLU:H	1.39	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:116:ASN:ND2	2:E:116:ASN:H	1.71	0.87
2:B:116:ASN:HD22	2:B:116:ASN:N	1.72	0.87
3:D:501:ANP:H8	3:D:501:ANP:H3'	1.57	0.87
1:D:393:GLU:OE2	1:D:394:GLY:N	2.07	0.86
2:E:116:ASN:N	2:E:116:ASN:HD22	1.75	0.84
2:E:116:ASN:ND2	2:E:116:ASN:N	2.23	0.83
1:D:232:LEU:HD11	1:D:243:ILE:HD12	1.62	0.79
1:D:361:PRO:HD2	2:E:42:ARG:NH1	1.98	0.78
1:A:391:ARG:NH2	2:B:46:ASP:OD2	2.16	0.78
1:D:392:ARG:HA	1:D:393:GLU:HB3	1.66	0.77
2:B:93:GLN:HG3	2:B:99:ILE:HD11	1.65	0.77
1:D:65:ALA:O	3:D:501:ANP:O2B	2.02	0.77
2:B:111:PRO:O	2:B:113:GLY:HA3	1.84	0.77
1:D:245:LEU:HD13	4:D:626:HOH:O	1.84	0.76
2:E:115:ARG:HD2	2:E:118:HIS:O	1.88	0.74
1:D:59:GLU:HB3	1:D:71:SER:HB3	1.69	0.73
1:A:297:TYR:HA	1:A:300:SER:HB3	1.71	0.73
1:A:97:LYS:NZ	1:A:101:ARG:HH22	1.86	0.72
1:A:309:TRP:HA	1:A:312:VAL:HG13	1.71	0.72
2:B:127:GLY:HA2	2:B:133:ARG:CB	2.20	0.71
1:A:145:HIS:CD2	2:B:128:PRO:HA	2.26	0.70
2:B:116:ASN:N	2:B:116:ASN:ND2	2.38	0.70
1:D:147:ILE:HD12	2:E:120:LEU:HD21	1.73	0.70
3:D:501:ANP:O1A	3:D:501:ANP:N3B	2.25	0.69
1:D:277:VAL:HG22	1:D:278:LEU:HD12	1.74	0.69
2:E:43:ASP:OD1	2:E:43:ASP:N	2.26	0.69
1:A:220:GLU:HG2	1:A:221:TYR:H	1.59	0.68
3:D:501:ANP:H3'	3:D:501:ANP:C8	2.24	0.68
2:B:33:VAL:HG12	2:B:34:HIS:H	1.59	0.68
1:A:96:ALA:HB1	1:A:377:ILE:HG21	1.77	0.67
1:A:355:ASP:OD2	4:A:612:HOH:O	2.12	0.66
2:B:116:ASN:ND2	2:B:116:ASN:H	1.92	0.66
2:B:29:VAL:HG11	2:B:51:VAL:HG11	1.77	0.66
1:A:308:PRO:HG2	1:A:311:THR:HG23	1.78	0.64
1:D:140:MET:O	3:D:501:ANP:H2	1.96	0.64
2:B:113:GLY:N	4:B:204:HOH:O	2.31	0.64
1:A:120:ILE:HG21	1:A:385:ILE:HG12	1.80	0.63
1:D:181:ARG:HG2	1:D:243:ILE:HD11	1.80	0.63
1:A:393:GLU:HG2	1:A:395:ILE:HB	1.81	0.63
2:E:40:LEU:HB2	2:E:43:ASP:OD1	1.98	0.63
2:B:76:GLU:O	2:B:79:LYS:N	2.28	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:391:ARG:O	1:D:393:GLU:HB3	1.99	0.62
2:B:34:HIS:HB2	2:B:35:SER:C	2.25	0.62
1:D:309:TRP:HA	1:D:312:VAL:HG13	1.82	0.62
1:A:203:MET:SD	1:A:221:TYR:HA	2.40	0.61
1:D:387:ASP:O	1:D:391:ARG:HB3	2.01	0.60
1:D:142:SER:OG	1:D:143:ASP:N	2.34	0.60
1:A:147:ILE:HG23	2:B:120:LEU:HD11	1.84	0.59
1:A:234:LEU:HD21	1:A:271:LEU:HD11	1.83	0.59
2:B:127:GLY:O	2:B:129:SER:N	2.36	0.59
2:E:23:ILE:HD12	2:E:24:PRO:HD2	1.85	0.59
2:B:23:ILE:HG13	2:B:26:SER:HB2	1.85	0.59
1:A:291:ALA:C	1:A:293:ARG:HA	2.28	0.58
1:D:206:GLY:HA2	1:D:217:PHE:HE1	1.67	0.58
1:D:228:ARG:HB3	1:D:233:MET:HE2	1.85	0.58
1:D:220:GLU:OE2	1:D:221:TYR:N	2.36	0.58
1:D:120:ILE:HG21	1:D:385:ILE:HG12	1.85	0.57
1:A:245:LEU:HD22	1:A:336:ILE:O	2.05	0.57
1:A:350:TYR:OH	4:A:605:HOH:O	2.18	0.56
1:A:398:GLN:HG3	1:D:237:HIS:CG	2.40	0.56
1:A:140:MET:O	3:A:501:ANP:H2	2.04	0.56
1:A:146:GLN:O	1:A:150:SER:OG	2.23	0.56
1:D:214:HIS:C	1:D:216:TYR:H	2.14	0.55
1:A:59:GLU:O	1:A:70:SER:OG	2.22	0.55
1:A:296:ALA:O	1:A:297:TYR:HB3	2.07	0.55
1:A:369:ARG:HA	1:A:372:LEU:HD12	1.88	0.55
1:A:218:MET:HE1	1:A:232:LEU:O	2.06	0.55
1:D:184:LYS:NZ	4:D:614:HOH:O	2.39	0.54
1:D:349:LYS:NZ	2:E:117:ILE:HD11	2.22	0.54
1:A:97:LYS:HZ2	1:A:101:ARG:HH22	1.53	0.54
2:E:112:PRO:HA	2:E:114:GLU:N	2.16	0.54
1:A:393:GLU:OE2	1:A:393:GLU:O	2.26	0.54
2:B:116:ASN:C	2:B:117:ILE:HG12	2.27	0.54
1:D:126:PRO:HG2	1:D:130:PHE:H	1.72	0.54
1:D:281:PRO:HD2	1:D:298:ILE:HD13	1.90	0.54
1:A:161:TYR:CE2	2:B:115:ARG:HG3	2.43	0.54
2:B:122:VAL:HG13	2:B:125:ARG:HD2	1.91	0.53
1:D:185:PRO:HD3	1:D:251:ILE:HG12	1.91	0.53
2:E:37:PRO:HB2	2:E:38:GLN:HG2	1.90	0.53
1:D:265:LYS:HB2	1:D:269:HIS:ND1	2.24	0.53
1:D:360:ALA:HB3	2:E:42:ARG:NH1	2.24	0.53
1:A:94:THR:HA	1:A:97:LYS:HE3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:GLU:OE1	1:A:160:ARG:NH1	2.38	0.52
1:D:83:ILE:HD13	1:D:136:VAL:HG22	1.90	0.52
1:A:291:ALA:O	1:A:293:ARG:HA	2.08	0.52
1:D:161:TYR:CZ	2:E:115:ARG:HB2	2.45	0.52
1:D:386:GLU:HA	1:D:389:HIS:HB2	1.91	0.52
2:B:40:LEU:HB2	2:B:43:ASP:OD2	2.09	0.52
1:D:231:GLU:OE1	1:D:335:ARG:NH2	2.38	0.52
2:B:64:ASP:HA	2:B:84:TYR:CZ	2.45	0.52
1:A:300:SER:OG	1:A:301:LEU:N	2.43	0.52
2:B:66:ASP:N	2:B:66:ASP:OD1	2.40	0.52
1:A:200:ASP:OD2	3:A:501:ANP:O1B	2.28	0.51
1:A:148:ILE:HD13	1:A:254:GLU:HG2	1.92	0.51
1:D:291:ALA:O	1:D:293:ARG:HA	2.10	0.51
1:A:261:LEU:HD13	1:A:313:TYR:OH	2.10	0.51
1:D:240:THR:O	1:D:243:ILE:HG23	2.10	0.51
1:D:393:GLU:OE2	1:D:393:GLU:C	2.54	0.51
1:A:221:TYR:CG	1:A:222:VAL:N	2.79	0.50
1:D:166:LEU:HD11	1:D:248:VAL:HG22	1.93	0.50
2:E:34:HIS:HB2	2:E:35:SER:HA	1.92	0.50
1:A:240:THR:N	4:A:604:HOH:O	2.26	0.50
1:A:257:ALA:HB2	1:A:313:TYR:HE1	1.77	0.50
1:D:393:GLU:O	1:D:395:ILE:N	2.45	0.50
2:B:33:VAL:HG12	2:B:34:HIS:N	2.26	0.50
1:D:225:ARG:HB3	1:D:233:MET:HE1	1.93	0.50
1:D:395:ILE:HG23	2:E:31:TRP:CD1	2.47	0.49
1:A:357:PRO:HG3	2:B:58:THR:HG21	1.93	0.49
1:D:211:PRO:HB2	2:B:26:SER:HB3	1.94	0.49
1:D:293:ARG:CB	1:D:296:ALA:H	2.26	0.49
1:D:379:GLU:HA	1:D:382:VAL:HB	1.94	0.49
1:A:221:TYR:OH	1:A:228:ARG:HD2	2.12	0.49
2:B:115:ARG:HD2	2:B:118:HIS:N	2.27	0.49
1:D:392:ARG:HA	1:D:393:GLU:CB	2.40	0.48
1:D:97:LYS:O	1:D:99:THR:N	2.47	0.48
3:D:501:ANP:O2A	3:D:501:ANP:O1B	2.32	0.48
2:B:61:GLU:HG3	2:B:105:PHE:HB2	1.96	0.48
1:A:93:VAL:O	1:A:97:LYS:HG3	2.14	0.47
1:A:75:ARG:HH11	1:D:339:ALA:HB2	1.80	0.47
1:A:149:HIS:HB3	2:B:129:SER:CB	2.44	0.47
2:B:48:ILE:HD13	2:B:106:PRO:HG3	1.96	0.47
1:A:152:GLN:HG2	4:A:609:HOH:O	2.13	0.47
2:E:47:VAL:O	2:E:51:VAL:HG13	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:HIS:CE1	1:A:168:ARG:NH1	2.83	0.47
2:B:126:ALA:HB1	2:B:135:HIS:HB2	1.96	0.46
1:A:152:GLN:HG2	1:A:153:PRO:HD2	1.97	0.46
1:A:212:ALA:HA	1:A:213:GLU:HA	1.53	0.46
1:A:220:GLU:HG2	1:A:221:TYR:N	2.28	0.46
2:B:125:ARG:HH11	2:B:125:ARG:HG2	1.81	0.46
1:A:152:GLN:OE1	2:B:120:LEU:HD13	2.15	0.46
1:D:379:GLU:O	1:D:383:ALA:N	2.48	0.46
2:E:61:GLU:OE2	2:E:107:ARG:NH1	2.48	0.46
2:E:120:LEU:HD22	2:E:121:LYS:N	2.30	0.46
1:A:47:VAL:HB	1:A:68:VAL:HG11	1.97	0.46
1:A:146:GLN:OE1	2:B:125:ARG:HB3	2.16	0.46
1:A:275:MET:HE3	1:A:280:THR:HG22	1.98	0.46
1:A:120:ILE:O	1:A:389:HIS:HE1	1.99	0.46
1:A:161:TYR:CZ	2:B:115:ARG:HB2	2.51	0.46
2:B:43:ASP:O	2:B:47:VAL:HG23	2.15	0.46
1:A:142:SER:HA	2:B:125:ARG:NH1	2.31	0.46
1:A:392:ARG:HA	1:A:393:GLU:HB3	1.98	0.46
2:B:134:HIS:O	2:B:135:HIS:CG	2.69	0.46
1:A:141:GLU:OE1	1:A:192:GLU:N	2.40	0.45
1:A:58:ILE:O	1:A:59:GLU:HG2	2.17	0.45
1:A:237:HIS:CG	1:A:237:HIS:O	2.69	0.45
1:D:352:ASP:HB3	1:D:355:ASP:HB3	1.98	0.45
2:B:41:PHE:HB2	2:B:78:MET:HG3	1.99	0.45
2:B:132:GLY:HA3	2:B:133:ARG:HA	1.44	0.45
1:D:77:THR:HG22	1:D:79:GLN:H	1.82	0.45
1:D:296:ALA:HA	1:D:299:GLN:HB2	1.98	0.45
2:E:110:LYS:HD2	2:E:114:GLU:HB3	1.99	0.45
1:A:352:ASP:HB3	1:A:355:ASP:HB3	1.98	0.45
1:D:176:ALA:HB2	1:D:359:CYS:HB3	1.98	0.45
1:D:392:ARG:H	1:D:392:ARG:HG3	1.48	0.45
1:D:95:ASN:N	1:D:95:ASN:OD1	2.50	0.45
1:A:145:HIS:HD2	2:B:128:PRO:HA	1.80	0.45
2:E:103:GLN:HG2	2:E:105:PHE:CZ	2.52	0.44
1:D:234:LEU:HD21	1:D:271:LEU:HD11	2.00	0.44
1:D:218:MET:HE1	1:D:232:LEU:HB3	1.99	0.44
1:A:111:HIS:HD2	1:A:172:TYR:CD2	2.35	0.44
2:B:116:ASN:HD22	2:B:116:ASN:H	1.44	0.44
1:D:82:ALA:HB2	1:D:139:LEU:HD23	1.99	0.44
2:B:42:ARG:HD3	2:B:42:ARG:HA	1.65	0.44
2:B:136:HIS:O	2:B:137:HIS:HB2	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:395:ILE:HG23	2:B:31:TRP:CD1	2.53	0.44
2:B:93:GLN:C	2:B:96:GLY:H	2.25	0.44
3:D:501:ANP:C8	3:D:501:ANP:C3'	2.86	0.43
2:B:21:ILE:HG12	2:B:104:ILE:HD11	2.00	0.43
1:D:209:THR:O	1:D:214:HIS:HE1	2.01	0.43
1:A:384:GLU:C	1:A:386:GLU:H	2.27	0.43
1:D:143:ASP:O	1:D:147:ILE:HG12	2.18	0.43
1:D:254:GLU:O	1:D:258:ARG:N	2.52	0.43
2:E:116:ASN:C	2:E:117:ILE:HG12	2.43	0.43
1:A:58:ILE:HG13	2:E:76:GLU:HG3	2.00	0.42
1:D:144:LEU:HB3	1:D:185:PRO:HB3	2.00	0.42
1:A:167:LEU:HD23	1:A:167:LEU:HA	1.80	0.42
1:D:72:ALA:HB3	1:D:81:VAL:HG12	2.01	0.42
1:D:69:VAL:HG12	1:D:70:SER:H	1.85	0.42
1:A:146:GLN:OE1	2:B:126:ALA:N	2.51	0.42
2:E:48:ILE:O	2:E:52:LEU:N	2.47	0.42
2:B:114:GLU:H	2:B:114:GLU:HG2	1.49	0.42
1:A:142:SER:HB2	2:B:122:VAL:HG13	2.01	0.42
1:D:98:ARG:NH2	3:D:501:ANP:O1G	2.53	0.42
1:D:122:ARG:HA	1:D:123:PRO:HD2	1.84	0.42
1:D:86:ILE:CG2	1:D:89:ALA:HB2	2.50	0.42
1:A:125:VAL:HA	1:A:126:PRO:HA	1.92	0.42
1:A:231:GLU:HA	1:A:236:LEU:HD12	2.02	0.41
1:A:310:GLU:O	1:A:314:PRO:HA	2.19	0.41
1:D:121:LEU:HD12	1:D:121:LEU:HA	1.71	0.41
1:A:233:MET:HE3	1:A:271:LEU:HD13	2.02	0.41
1:A:289:VAL:O	1:A:295:ARG:NH1	2.52	0.41
2:E:120:LEU:HD22	2:E:121:LYS:H	1.84	0.41
1:A:323:LEU:O	1:A:327:MET:HG3	2.20	0.41
2:E:17:LEU:HD12	2:E:18:VAL:H	1.85	0.41
2:B:122:VAL:CG1	2:B:125:ARG:HD2	2.51	0.41
1:D:98:ARG:HB3	1:D:99:THR:H	1.74	0.41
2:E:23:ILE:HD13	2:E:52:LEU:HD21	2.03	0.41
1:D:368:ASP:HB3	1:D:369:ARG:H	1.72	0.41
1:A:293:ARG:HG2	1:A:295:ARG:HE	1.86	0.41
2:E:84:TYR:CD1	2:E:84:TYR:C	2.99	0.41
2:B:86:TYR:HD1	2:B:86:TYR:HA	1.74	0.41
2:B:74:SER:C	2:B:75:ASP:O	2.58	0.40
2:E:64:ASP:CB	2:E:68:ASP:HB2	2.51	0.40
1:D:55:TYR:HD2	1:D:121:LEU:HD21	1.86	0.40
1:D:310:GLU:O	1:D:314:PRO:HA	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:98:LEU:H	2:B:98:LEU:HD23	1.87	0.40
1:A:320:ALA:HB2	1:A:346:PHE:CE1	2.57	0.40
3:A:501:ANP:H8	3:A:501:ANP:O5'	2.21	0.40
1:D:86:ILE:HG22	1:D:89:ALA:HB2	2.03	0.40
1:A:399:ILE:O	1:A:399:ILE:HG13	2.22	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	353/442 (80%)	324 (92%)	25 (7%)	4 (1%)	11	25
1	D	351/442 (79%)	303 (86%)	44 (12%)	4 (1%)	11	25
2	B	121/126 (96%)	103 (85%)	17 (14%)	1 (1%)	16	34
2	E	109/126 (86%)	101 (93%)	6 (6%)	2 (2%)	6	14
All	All	934/1136 (82%)	831 (89%)	92 (10%)	11 (1%)	10	23

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	222	VAL
1	A	385	ILE
1	D	367	PHE
2	E	16	VAL
1	D	385	ILE
1	A	216	TYR
1	A	372	LEU
1	D	368	ASP
1	D	371	ALA
2	B	117	ILE

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Mol	Chain	Res	Type
2	E	122	VAL

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	283/373 (76%)	246 (87%)	37 (13%)	4	8
1	D	262/373 (70%)	223 (85%)	39 (15%)	3	5
2	B	98/110 (89%)	77 (79%)	21 (21%)	1	2
2	E	92/110 (84%)	81 (88%)	11 (12%)	5	10
All	All	735/966 (76%)	627 (85%)	108 (15%)	3	6

All (108) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	VAL
1	A	48	THR
1	A	58	ILE
1	A	61	ILE
1	A	68	VAL
1	A	71	SER
1	A	97	LYS
1	A	124	THR
1	A	135	VAL
1	A	139	LEU
1	A	156	LEU
1	A	159	VAL
1	A	166	LEU
1	A	170	LEU
1	A	173	MET
1	A	196	LEU
1	A	203	MET
1	A	208	CYS
1	A	220	GLU
1	A	224	THR

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Mol	Chain	Res	Type
1	A	228	ARG
1	A	235	SER
1	A	245	LEU
1	A	268	VAL
1	A	285	VAL
1	A	305	GLN
1	A	311	THR
1	A	312	VAL
1	A	318	ARG
1	A	331	GLU
1	A	365	PHE
1	A	373	THR
1	A	377	ILE
1	A	392	ARG
1	A	393	GLU
1	A	398	GLN
1	A	399	ILE
1	D	59	GLU
1	D	63	ASN
1	D	69	VAL
1	D	81	VAL
1	D	83	ILE
1	D	90	PHE
1	D	92	VAL
1	D	95	ASN
1	D	121	LEU
1	D	141	GLU
1	D	156	LEU
1	D	159	VAL
1	D	166	LEU
1	D	170	LEU
1	D	196	LEU
1	D	219	THR
1	D	220	GLU
1	D	224	THR
1	D	228	ARG
1	D	243	ILE
1	D	245	LEU
1	D	265	LYS
1	D	271	LEU
1	D	275	MET
1	D	277	VAL

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Mol	Chain	Res	Type
1	D	285	VAL
1	D	294	VAL
1	D	298	ILE
1	D	299	GLN
1	D	311	THR
1	D	312	VAL
1	D	326	ARG
1	D	331	GLU
1	D	365	PHE
1	D	368	ASP
1	D	389	HIS
1	D	391	ARG
1	D	392	ARG
1	D	393	GLU
2	E	43	ASP
2	E	45	LEU
2	E	52	LEU
2	E	57	THR
2	E	63	GLU
2	E	65	GLU
2	E	94	VAL
2	E	98	LEU
2	E	116	ASN
2	E	117	ILE
2	E	120	LEU
2	B	17	LEU
2	B	26	SER
2	B	34	HIS
2	B	52	LEU
2	B	65	GLU
2	B	66	ASP
2	B	76	GLU
2	B	88	THR
2	B	89	VAL
2	B	93	GLN
2	B	94	VAL
2	B	107	ARG
2	B	114	GLU
2	B	115	ARG
2	B	116	ASN
2	B	117	ILE
2	B	120	LEU

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Mol	Chain	Res	Type
2	B	121	LYS
2	B	124	THR
2	B	130	GLN
2	B	137	HIS

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

Mol	Chain	Res	Type
1	A	108	HIS
1	A	149	HIS
1	D	88	ASN
1	D	145	HIS
2	E	116	ASN
2	B	93	GLN
2	B	116	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](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ANP	D	501	-	33,33,33	2.56	14 (42%)	45,52,52	2.65	14 (31%)
3	ANP	A	501	-	33,33,33	2.56	13 (39%)	45,52,52	2.53	15 (33%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ANP	D	501	-	-	3/18/38/38	0/3/3/3
3	ANP	A	501	-	-	3/18/38/38	0/3/3/3

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	501	ANP	PB-O1B	8.64	1.59	1.46
3	A	501	ANP	PB-O1B	8.27	1.58	1.46
3	A	501	ANP	C2'-C3'	-5.01	1.39	1.53
3	D	501	ANP	C2'-C3'	-4.63	1.40	1.53
3	A	501	ANP	PB-N3B	4.58	1.75	1.63
3	D	501	ANP	PB-N3B	4.53	1.75	1.63
3	D	501	ANP	PA-O3A	-4.12	1.55	1.59
3	A	501	ANP	O4'-C4'	-3.77	1.36	1.45
3	A	501	ANP	C6-N6	3.59	1.43	1.34
3	D	501	ANP	C6-N6	3.52	1.43	1.34
3	A	501	ANP	PA-O3A	-3.37	1.55	1.59
3	D	501	ANP	O4'-C4'	-3.37	1.37	1.45
3	A	501	ANP	C5-N7	-2.76	1.34	1.39
3	D	501	ANP	C3'-C4'	-2.64	1.46	1.53
3	A	501	ANP	C3'-C4'	-2.58	1.46	1.53
3	A	501	ANP	O2'-C2'	-2.57	1.36	1.43
3	A	501	ANP	O3'-C3'	-2.57	1.36	1.43
3	A	501	ANP	C2'-C1'	-2.56	1.45	1.53
3	A	501	ANP	PG-O1G	2.53	1.50	1.46
3	D	501	ANP	C2'-C1'	-2.45	1.45	1.53
3	A	501	ANP	C8-N9	-2.41	1.33	1.37
3	D	501	ANP	O3'-C3'	-2.39	1.37	1.43
3	D	501	ANP	C5-N7	-2.35	1.34	1.39
3	D	501	ANP	PG-O1G	2.35	1.49	1.46
3	D	501	ANP	O2'-C2'	-2.23	1.37	1.43
3	D	501	ANP	PB-O3A	-2.08	1.56	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	501	ANP	C8-N9	-2.03	1.34	1.37

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	501	ANP	O1B-PB-N3B	-10.26	96.66	111.77
3	A	501	ANP	O1B-PB-N3B	-9.58	97.66	111.77
3	A	501	ANP	C5-C4-N3	-5.53	119.10	126.72
3	D	501	ANP	C5-C4-N3	-5.37	119.32	126.72
3	D	501	ANP	O2B-PB-O3A	5.01	121.35	104.64
3	A	501	ANP	O2B-PB-O3A	4.91	121.03	104.64
3	D	501	ANP	O2B-PB-O1B	4.57	119.68	109.87
3	D	501	ANP	N3-C4-N9	4.21	134.32	127.17
3	A	501	ANP	N3-C4-N9	4.20	134.31	127.17
3	D	501	ANP	N9-C8-N7	-4.12	108.08	113.94
3	A	501	ANP	N9-C8-N7	-4.10	108.12	113.94
3	A	501	ANP	C5-N7-C8	3.83	109.47	103.45
3	D	501	ANP	C5-N7-C8	3.73	109.32	103.45
3	A	501	ANP	C4-C5-N7	-3.40	106.69	110.58
3	D	501	ANP	C4-N9-C8	3.39	109.29	105.74
3	D	501	ANP	N3-C2-N1	-3.26	123.65	128.58
3	A	501	ANP	C4-N9-C8	3.22	109.12	105.74
3	D	501	ANP	C4-C5-N7	-3.19	106.94	110.58
3	A	501	ANP	O2B-PB-O1B	3.06	116.42	109.87
3	D	501	ANP	C2-N3-C4	2.85	118.78	111.83
3	A	501	ANP	N3-C2-N1	-2.84	124.28	128.58
3	A	501	ANP	C2-N3-C4	2.77	118.59	111.83
3	D	501	ANP	O3A-PB-N3B	-2.68	99.16	106.59
3	A	501	ANP	O5'-C5'-C4'	2.62	117.91	108.99
3	D	501	ANP	C2'-C1'-N9	-2.48	107.14	113.30
3	A	501	ANP	C2'-C1'-N9	-2.47	107.17	113.30
3	D	501	ANP	O5'-C5'-C4'	2.40	117.16	108.99
3	A	501	ANP	C6-C5-C4	2.16	120.13	117.18
3	A	501	ANP	O2G-PG-O3G	-2.10	101.95	107.59

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	ANP	PB-N3B-PG-O1G
3	A	501	ANP	PA-O3A-PB-O1B
3	A	501	ANP	PA-O3A-PB-O2B

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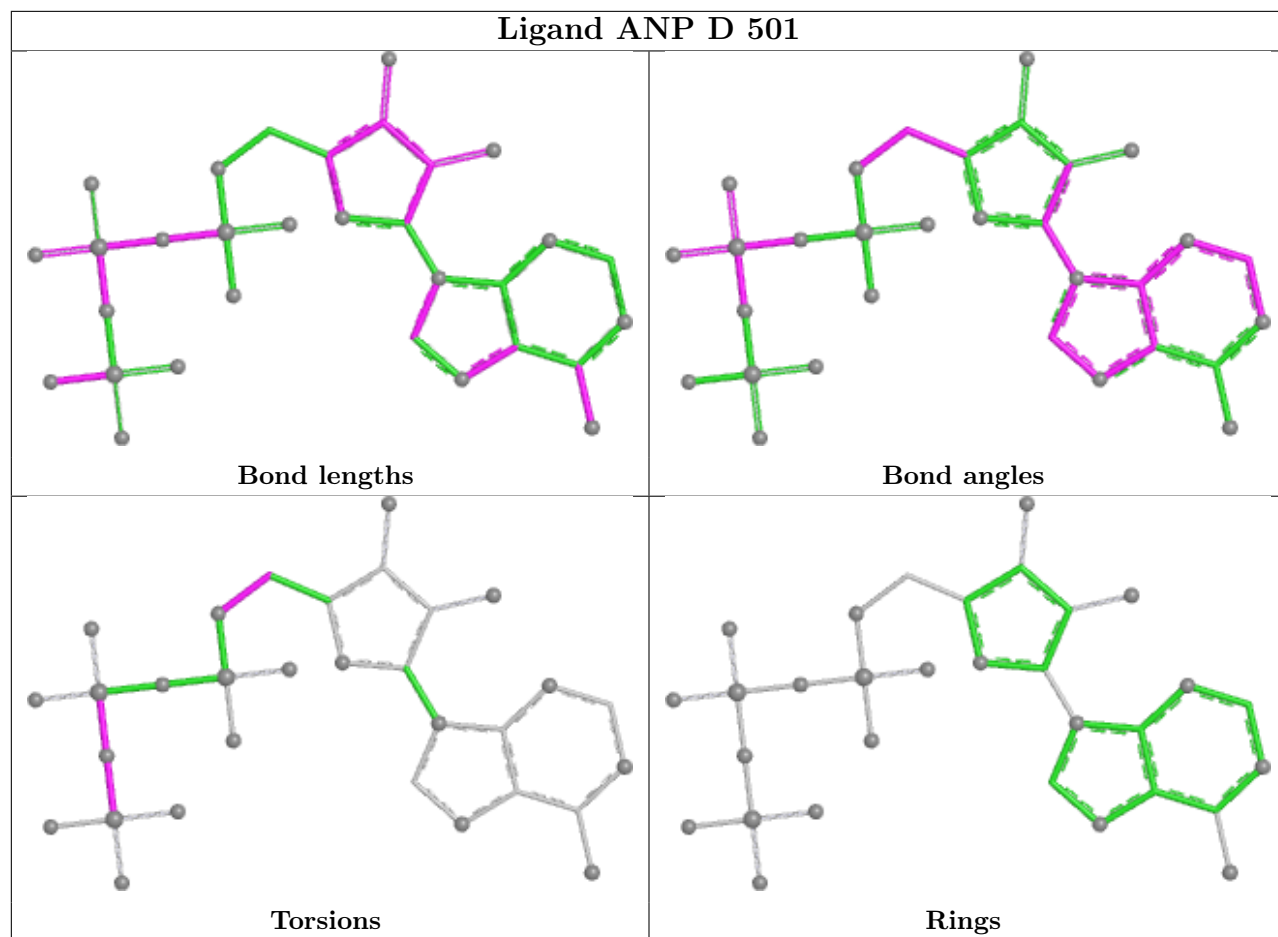
Mol	Chain	Res	Type	Atoms
3	D	501	ANP	PB-N3B-PG-O1G
3	D	501	ANP	PG-N3B-PB-O1B
3	D	501	ANP	C4'-C5'-O5'-PA

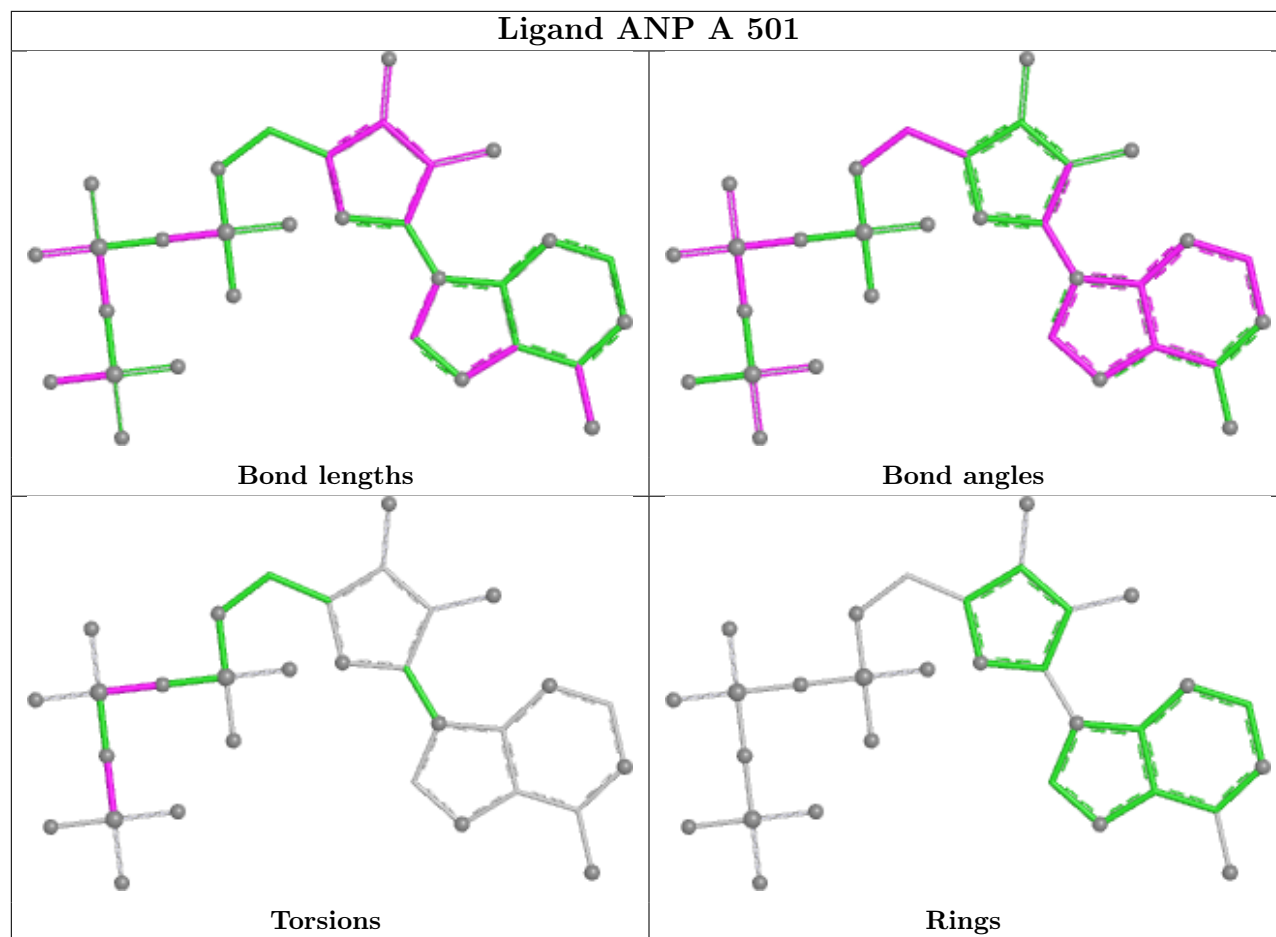
There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	501	ANP	8	0
3	A	501	ANP	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.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	355/442 (80%)	0.20	10 (2%) 55 49	49, 82, 141, 206	0
1	D	353/442 (79%)	0.29	8 (2%) 61 55	49, 84, 162, 184	0
2	B	123/126 (97%)	0.49	5 (4%) 41 36	60, 92, 131, 152	0
2	E	111/126 (88%)	0.38	5 (4%) 38 32	62, 89, 150, 182	0
All	All	942/1136 (82%)	0.29	28 (2%) 52 47	49, 85, 153, 206	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	372	LEU	3.6
2	E	14	GLY	3.4
2	E	123	ASN	3.2
2	B	128	PRO	3.0
1	D	120	ILE	2.9
1	A	371	ALA	2.9
2	B	136	HIS	2.9
2	E	124	THR	2.8
1	A	399	ILE	2.7
2	E	122	VAL	2.7
2	E	15	SER	2.7
2	B	138	HIS	2.6
1	A	291	ALA	2.4
1	A	395	ILE	2.4
1	A	222	VAL	2.4
1	D	125	VAL	2.3
1	D	360	ALA	2.3
1	D	134	TYR	2.3
1	A	125	VAL	2.3
1	A	215	GLN	2.2
1	D	124	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	96	ALA	2.2
1	D	93	VAL	2.2
1	D	399	ILE	2.2
1	A	301	LEU	2.1
1	A	221	TYR	2.1
2	B	135	HIS	2.1
2	B	35	SER	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

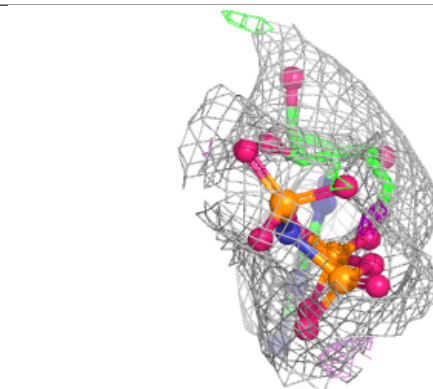
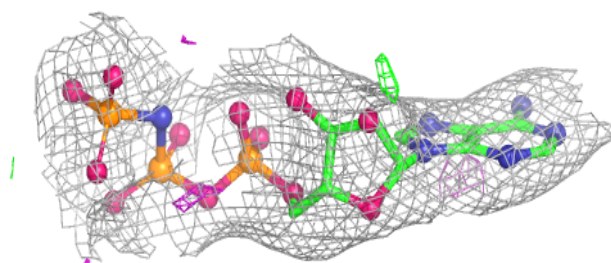
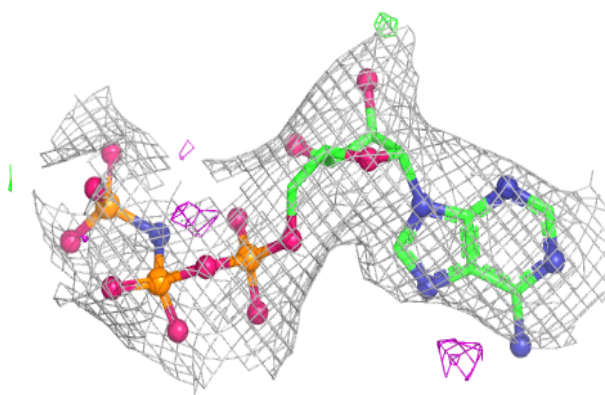
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ANP	D	501	31/31	0.80	0.12	90,121,213,278	0
3	ANP	A	501	31/31	0.92	0.11	66,80,236,380	0

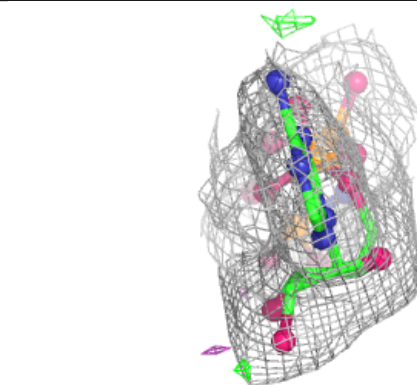
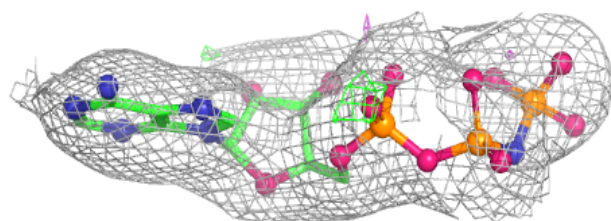
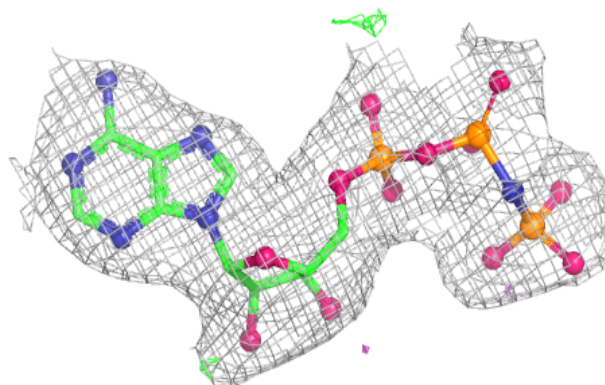
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ANP 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)

**Electron density around ANP A 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.