



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:10 PM UTC

PDB ID : 1KI3 / pdb_00001ki3
Title : CRYSTAL STRUCTURE OF THYMIDINE KINASE FROM HERPES SIMPLEX VIRUS TYPE I COMPLEXED WITH PENCICLOVIR
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Deposited on : 1998-05-15
Resolution : 2.37 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

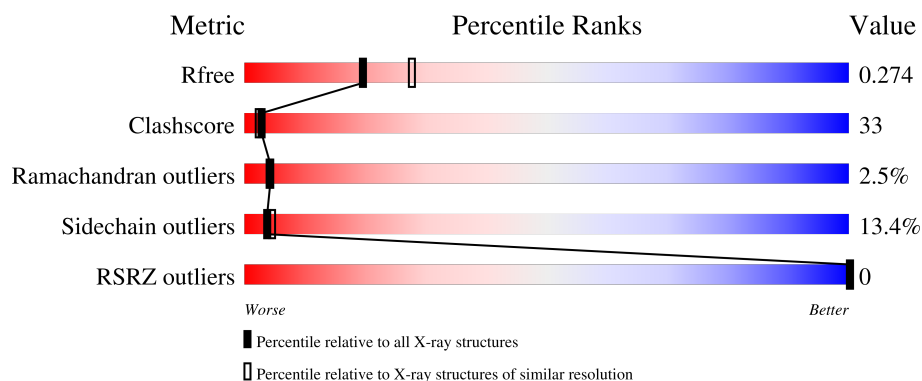
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.37 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	331	
1	B	331	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4869 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 THYMIDINE KINASE.

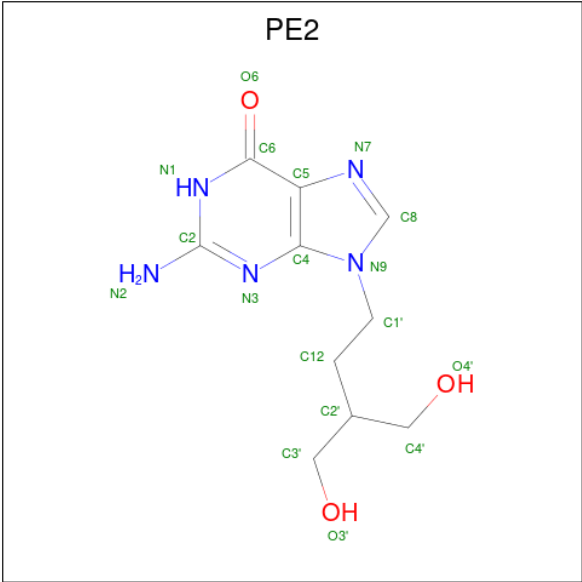
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	0	3
			2339	1490	411	423	15			
1	B	311	Total	C	N	O	S	0	0	3
			2352	1500	413	424	15			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 9-(4-HYDROXY-3-(HYDROXYMETHYL)BUT-1-YL)GUANINE (CCD ID: PE2) (formula: C₁₀H₁₅N₅O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			18	10	5	3		
3	B	1	Total	C	N	O	0	0
			18	10	5	3		

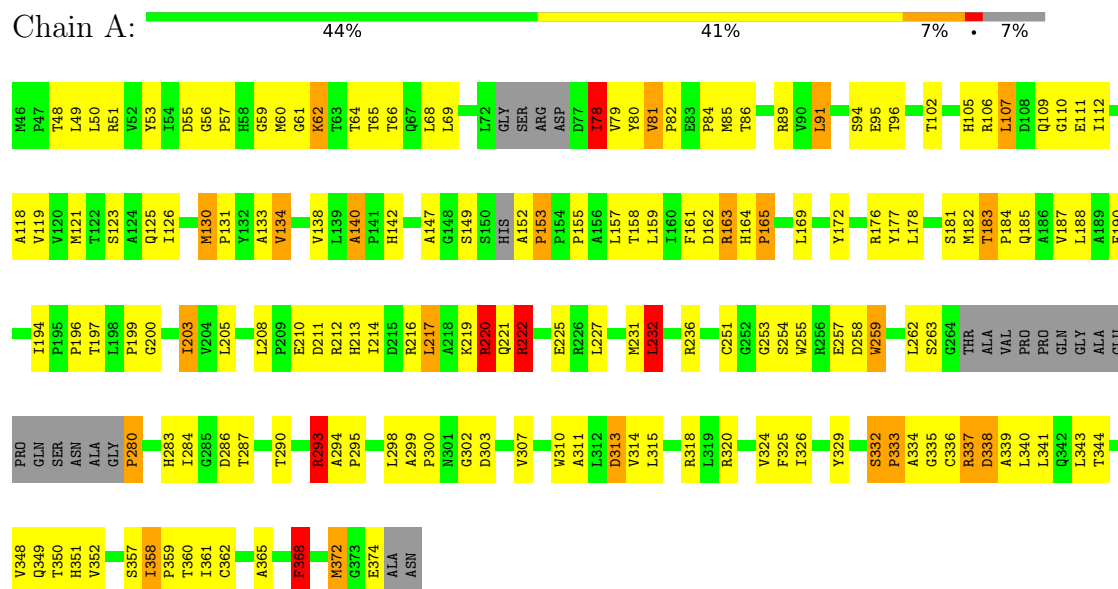
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	69	Total	O	0	0
			69	69		
4	B	63	Total	O	0	0
			63	63		

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: THYMIDINE KINASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	113.80Å 117.70Å 108.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 2.37 12.00 – 2.37	Depositor EDS
% Data completeness (in resolution range)	70.4 (12.00-2.37) 70.4 (12.00-2.37)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.35Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.200 , 0.314 0.235 , 0.274	Depositor DCC
R_{free} test set	2096 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.7	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.057 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4869	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, PE2

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	0.94	2/2392 (0.1%)	1.24	25/3267 (0.8%)
1	B	0.98	4/2406 (0.2%)	1.20	14/3287 (0.4%)
All	All	0.96	6/4798 (0.1%)	1.22	39/6554 (0.6%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	149	SER	C-N	-5.51	1.25	1.33
1	B	316	ALA	C-O	-5.26	1.18	1.24
1	B	337	ARG	C-O	-5.25	1.18	1.24
1	B	100	ILE	CA-CB	5.19	1.60	1.54
1	A	263	SER	C-N	-5.07	1.26	1.33
1	B	306	ASN	C-O	-5.06	1.18	1.24

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	293	ARG	N-CA-C	-10.93	100.22	112.57
1	A	91	LEU	N-CA-C	9.13	122.06	111.11
1	A	81	VAL	CA-C-N	7.96	129.79	119.84
1	A	81	VAL	C-N-CA	7.96	129.79	119.84
1	B	222	ARG	N-CA-C	7.55	119.42	109.83
1	A	368	PHE	N-CA-C	-7.49	103.05	111.14
1	B	81	VAL	CA-C-N	7.46	127.37	120.21
1	B	81	VAL	C-N-CA	7.46	127.37	120.21
1	A	280	PRO	N-CA-CB	7.30	111.03	103.00
1	B	205	LEU	CA-C-N	-6.91	116.95	121.65
1	B	205	LEU	C-N-CA	-6.91	116.95	121.65
1	B	354	THR	N-CA-C	-6.84	101.50	110.39
1	B	67	GLN	N-CA-C	-6.24	104.17	110.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	351	HIS	N-CA-C	-6.24	101.29	110.52
1	A	360	THR	N-CA-C	-6.15	104.57	111.28
1	A	222	ARG	N-CA-C	6.10	123.30	109.81
1	A	253	GLY	N-CA-C	5.82	119.45	111.54
1	B	228	ASP	N-CA-C	-5.72	97.87	107.99
1	A	332	SER	CA-C-N	5.67	126.93	119.84
1	A	332	SER	C-N-CA	5.67	126.93	119.84
1	B	312	LEU	N-CA-C	-5.66	105.20	111.71
1	A	183	THR	N-CA-C	5.59	116.74	109.64
1	A	358	ILE	CA-C-N	5.45	125.10	119.05
1	A	358	ILE	C-N-CA	5.45	125.10	119.05
1	A	231	MET	N-CA-C	-5.44	105.91	112.54
1	A	313	ASP	N-CA-C	-5.40	105.40	111.28
1	A	220	ARG	N-CA-C	-5.39	99.32	110.80
1	B	320	ARG	N-CA-C	5.30	118.54	111.75
1	A	94	SER	N-CA-C	5.25	117.29	108.73
1	A	336	CYS	N-CA-C	-5.24	105.16	112.45
1	B	292	PHE	N-CA-C	5.21	119.34	112.25
1	B	222	ARG	CA-C-N	5.21	126.36	119.84
1	B	222	ARG	C-N-CA	5.21	126.36	119.84
1	A	78	ILE	N-CA-C	5.20	115.33	107.80
1	A	232	LEU	N-CA-C	-5.19	105.31	111.69
1	B	104	GLN	N-CA-C	-5.05	105.77	111.28
1	A	107	LEU	N-CA-C	-5.05	105.47	110.97
1	A	140	ALA	N-CA-C	5.03	120.94	109.81
1	A	326	ILE	N-CA-C	5.02	115.91	108.23

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	2339	0	2349	155	0
1	B	2352	0	2368	157	0
2	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	5	0	0	0	0
3	A	18	0	15	0	0
3	B	18	0	15	5	0
4	A	69	0	0	6	0
4	B	63	0	0	4	0
All	All	4869	0	4747	310	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 33.

All (310) 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:254:SER:HB3	1:A:257:GLU:HG2	1.30	1.10
1:A:84:PRO:HG3	1:A:368:PHE:HZ	1.19	1.02
1:B:62:LYS:HE2	1:B:163:ARG:HH12	1.37	0.89
1:A:232:LEU:HD13	1:A:236:ARG:HD2	1.54	0.88
1:A:84:PRO:HG3	1:A:368:PHE:CZ	2.09	0.85
1:A:334:ALA:HA	1:A:337:ARG:HB3	1.58	0.85
1:A:130:MET:HB3	1:A:131:PRO:HD3	1.59	0.85
1:A:85:MET:HE2	1:A:89:ARG:HG2	1.58	0.84
1:A:251:CYS:SG	1:A:280:PRO:O	2.37	0.82
1:A:78:ILE:HG12	1:A:158:THR:HB	1.64	0.80
1:A:57:PRO:HD2	1:A:60:MET:HE2	1.61	0.80
1:B:327:LEU:HD22	1:B:340:LEU:HB2	1.64	0.80
1:A:121:MET:HG3	1:A:182:MET:HE2	1.63	0.79
1:B:211:ASP:O	1:B:215:ASP:HB2	1.85	0.77
1:B:358:ILE:HB	1:B:359:PRO:HD3	1.65	0.77
1:B:63:THR:O	1:B:67:GLN:HG2	1.85	0.77
1:B:88:TRP:CE3	1:B:97:ILE:HD11	2.20	0.77
1:B:178:LEU:HD11	1:B:291:LEU:HD22	1.66	0.76
1:B:149:SER:HA	1:B:347:MET:O	1.86	0.76
3:B:2:PE2:H1'2	4:B:527:HOH:O	1.84	0.76
1:A:61:GLY:O	1:A:65:THR:HG23	1.86	0.76
1:A:68:LEU:HB3	1:A:337:ARG:HG3	1.68	0.75
1:B:183:THR:O	1:B:187:VAL:HG23	1.86	0.73
1:B:216:ARG:O	1:B:219:LYS:HG2	1.89	0.73
1:A:217:LEU:HD12	1:A:232:LEU:HG	1.70	0.73
1:B:208:LEU:HD11	1:B:212:ARG:HD2	1.69	0.72
1:A:126:ILE:HD11	1:B:126:ILE:HD11	1.72	0.71
1:B:210:GLU:HG3	1:B:214:ILE:HD13	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:LEU:HD22	1:A:341:LEU:HG	1.72	0.70
1:A:184:PRO:HG2	1:A:185:GLN:OE1	1.91	0.70
1:B:246:VAL:O	1:B:250:GLN:HG3	1.92	0.70
1:A:232:LEU:O	1:A:236:ARG:HG3	1.92	0.70
1:B:248:TYR:HB2	1:B:282:PRO:HG2	1.74	0.70
1:A:314:VAL:HG13	1:A:318:ARG:NH1	2.07	0.69
1:A:85:MET:HB3	1:A:374:GLU:HB3	1.74	0.69
1:A:164:HIS:CG	1:A:165:PRO:HD2	2.27	0.69
1:A:85:MET:HE2	1:A:89:ARG:CG	2.22	0.68
1:B:121:MET:HB3	1:B:182:MET:HE2	1.76	0.68
1:B:69:LEU:HD11	1:B:341:LEU:HD23	1.74	0.68
1:B:258:ASP:HB2	1:B:284:ILE:HD12	1.73	0.68
1:A:213:HIS:CE1	1:A:232:LEU:HD21	2.29	0.68
1:B:163:ARG:HH21	1:B:172:TYR:HE2	1.39	0.68
1:B:107:LEU:HA	1:B:112:ILE:HG22	1.76	0.68
1:A:49:LEU:HD12	1:A:157:LEU:O	1.94	0.67
1:A:130:MET:O	1:A:134:VAL:HG23	1.95	0.67
1:B:69:LEU:CD1	1:B:341:LEU:HD23	2.25	0.67
1:B:62:LYS:HE2	1:B:163:ARG:NH1	2.09	0.66
1:B:118:ALA:HA	1:B:181:SER:O	1.96	0.66
1:A:69:LEU:CD2	1:A:341:LEU:HG	2.26	0.66
1:B:121:MET:HG3	1:B:181:SER:HB2	1.78	0.66
1:A:142:HIS:HB3	1:A:361:ILE:HD11	1.78	0.65
1:B:53:TYR:HB2	1:B:203:ILE:HD13	1.78	0.65
1:A:211:ASP:HA	1:A:214:ILE:HD12	1.77	0.65
1:A:358:ILE:HB	1:A:359:PRO:HD3	1.79	0.65
1:A:118:ALA:HA	1:A:181:SER:O	1.96	0.64
1:B:85:MET:HE2	1:B:89:ARG:HG2	1.79	0.64
1:B:85:MET:HE3	1:B:88:TRP:HE3	1.63	0.64
1:B:226:ARG:HA	4:B:557:HOH:O	1.98	0.63
1:A:79:VAL:HG12	1:A:80:TYR:N	2.14	0.62
1:B:49:LEU:CD1	1:B:157:LEU:HD23	2.30	0.61
1:A:212:ARG:NH2	1:A:216:ARG:HE	1.97	0.61
1:B:62:LYS:CE	1:B:163:ARG:HH12	2.12	0.61
1:A:57:PRO:HD2	1:A:60:MET:CE	2.30	0.61
1:B:112:ILE:HG12	1:B:116:ASP:HB2	1.82	0.60
1:A:176:ARG:HB2	1:A:182:MET:HE3	1.84	0.60
1:B:116:ASP:O	1:B:120:VAL:HG23	2.01	0.60
1:B:130:MET:O	1:B:134:VAL:HG23	2.01	0.60
1:B:212:ARG:O	1:B:216:ARG:HG3	2.01	0.60
1:B:262:LEU:CD1	1:B:284:ILE:HG23	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:ALA:O	1:A:152:ALA:N	2.35	0.59
1:B:97:ILE:HG12	1:B:128:MET:HE1	1.84	0.59
1:B:373:GLY:O	1:B:374:GLU:HB2	2.01	0.59
1:B:298:LEU:HD22	1:B:302:GLY:O	2.02	0.59
1:A:79:VAL:HG12	1:A:80:TYR:H	1.67	0.59
1:A:155:PRO:HD3	1:A:349:GLN:NE2	2.17	0.59
1:A:85:MET:CE	1:A:89:ARG:HG2	2.33	0.58
1:B:134:VAL:O	1:B:138:VAL:HG22	2.03	0.58
1:A:109:GLN:O	1:A:111:GLU:HG3	2.04	0.58
3:B:2:PE2:O3'	3:B:2:PE2:H8	2.03	0.58
1:A:155:PRO:HD3	1:A:349:GLN:HE22	1.67	0.58
1:B:146:GLU:CD	1:B:348:VAL:HG11	2.28	0.58
1:B:215:ASP:O	1:B:218:ALA:HB3	2.04	0.58
1:A:334:ALA:O	1:A:338:ASP:N	2.34	0.58
1:B:208:LEU:HD12	1:B:209:PRO:HD2	1.85	0.58
1:B:214:ILE:HG13	1:B:229:LEU:HD22	1.85	0.58
1:B:327:LEU:CD2	1:B:340:LEU:HB2	2.31	0.58
1:B:292:PHE:CD2	1:B:312:LEU:HD21	2.40	0.57
1:A:62:LYS:NZ	1:A:62:LYS:HB2	2.20	0.57
1:A:318:ARG:NH1	4:A:585:HOH:O	2.36	0.57
1:B:50:LEU:HB2	1:B:349:GLN:HB3	1.84	0.57
1:A:59:GLY:O	1:A:213:HIS:HD2	1.87	0.57
1:B:49:LEU:HD13	1:B:157:LEU:HD23	1.87	0.57
1:A:232:LEU:CD1	1:A:236:ARG:HD2	2.32	0.56
1:A:222:ARG:HB2	1:A:225:GLU:HG3	1.88	0.56
1:B:246:VAL:HG21	1:B:324:VAL:HG21	1.87	0.56
1:B:139:LEU:HD11	1:B:143:ILE:HD11	1.87	0.56
1:A:134:VAL:O	1:A:138:VAL:HG23	2.05	0.56
1:A:130:MET:SD	1:A:134:VAL:CG2	2.94	0.55
1:A:163:ARG:NH2	1:A:172:TYR:OH	2.39	0.55
1:B:130:MET:HB3	1:B:131:PRO:HD3	1.89	0.55
1:A:258:ASP:HB2	1:A:284:ILE:HD12	1.89	0.55
1:A:333:PRO:C	1:A:335:GLY:H	2.14	0.55
1:A:69:LEU:HD23	1:A:337:ARG:HD3	1.88	0.55
1:A:287:THR:O	1:A:290:THR:HG22	2.07	0.55
1:A:60:MET:HG2	1:A:213:HIS:CD2	2.42	0.55
1:A:205:LEU:HD12	1:A:324:VAL:HG21	1.88	0.55
1:B:222:ARG:O	1:B:225:GLU:HB3	2.06	0.54
1:A:219:LYS:HG2	1:A:220:ARG:N	2.23	0.54
1:B:213:HIS:HB2	1:B:236:ARG:HH21	1.72	0.54
1:A:69:LEU:HD11	1:A:340:LEU:HG	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:THR:HG22	1:A:91:LEU:HG	1.90	0.54
1:A:208:LEU:O	1:A:236:ARG:NH2	2.41	0.54
1:A:185:GLN:OE1	1:A:185:GLN:N	2.39	0.54
1:B:69:LEU:HD11	1:B:341:LEU:CD2	2.37	0.54
1:B:142:HIS:HA	1:B:353:THR:HB	1.90	0.54
1:A:121:MET:HG3	1:A:182:MET:CE	2.35	0.53
1:B:69:LEU:HD22	1:B:340:LEU:CD2	2.39	0.53
1:B:362:CYS:O	1:B:366:ARG:HG3	2.08	0.53
1:B:85:MET:HE3	1:B:88:TRP:CE3	2.42	0.53
1:B:247:ARG:O	1:B:248:TYR:C	2.51	0.53
1:B:262:LEU:HD11	1:B:284:ILE:HG23	1.90	0.53
1:B:142:HIS:CD2	1:B:357:SER:OG	2.62	0.53
1:B:88:TRP:HE3	1:B:97:ILE:HD11	1.74	0.53
1:B:222:ARG:HD2	1:B:222:ARG:N	2.22	0.53
1:B:121:MET:CE	1:B:125:GLN:HE21	2.21	0.53
1:B:142:HIS:HD2	1:B:357:SER:OG	1.91	0.53
1:B:314:VAL:O	1:B:318:ARG:HG2	2.09	0.53
1:A:220:ARG:HH11	1:A:220:ARG:HB3	1.74	0.53
1:A:298:LEU:HA	1:A:303:ASP:O	2.08	0.52
1:A:123:SER:O	1:A:126:ILE:HB	2.09	0.52
1:B:244:ASN:HA	1:B:247:ARG:HE	1.75	0.52
1:A:130:MET:SD	1:A:134:VAL:HG21	2.50	0.52
1:B:146:GLU:OE1	1:B:348:VAL:HG11	2.09	0.52
1:B:214:ILE:HD12	1:B:232:LEU:CD1	2.39	0.52
1:A:102:THR:O	1:A:105:HIS:HB3	2.10	0.52
1:A:220:ARG:O	1:A:221:GLN:HG2	2.10	0.52
1:B:83:GLU:HG2	1:B:88:TRP:CH2	2.44	0.52
1:A:59:GLY:O	1:A:213:HIS:CD2	2.63	0.52
1:A:106:ARG:NH2	4:A:566:HOH:O	2.42	0.52
1:A:126:ILE:HD13	1:A:190:PHE:HE1	1.75	0.51
1:A:130:MET:HB3	1:A:131:PRO:CD	2.37	0.51
1:A:157:LEU:HD22	1:A:358:ILE:HG23	1.91	0.51
1:B:255:TRP:HA	1:B:284:ILE:HD12	1.92	0.51
1:B:289:PHE:HD1	1:B:312:LEU:HD22	1.75	0.51
1:B:220:ARG:HG2	1:B:221:GLN:N	2.25	0.51
1:B:232:LEU:O	1:B:236:ARG:HG3	2.10	0.51
1:A:203:ILE:O	1:A:324:VAL:HA	2.10	0.51
1:B:248:TYR:CE2	1:B:253:GLY:HA3	2.46	0.51
1:A:169:LEU:HD13	1:A:194:ILE:HA	1.92	0.51
1:A:106:ARG:NH1	1:A:112:ILE:HD13	2.26	0.50
1:A:78:ILE:HG12	1:A:158:THR:CB	2.39	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:LEU:HD12	1:A:324:VAL:CG2	2.41	0.50
1:A:333:PRO:O	1:A:334:ALA:HB3	2.11	0.50
1:A:56:GLY:O	1:A:163:ARG:NH2	2.44	0.50
1:B:219:LYS:O	1:B:220:ARG:HB2	2.11	0.50
1:A:183:THR:HG22	1:A:185:GLN:H	1.76	0.50
1:B:254:SER:HA	4:B:550:HOH:O	2.12	0.50
1:A:352:VAL:HB	1:A:357:SER:CB	2.42	0.50
1:A:51:ARG:NH1	1:A:53:TYR:OH	2.45	0.49
1:A:333:PRO:C	1:A:335:GLY:N	2.70	0.49
1:A:334:ALA:CA	1:A:337:ARG:HB3	2.39	0.49
1:B:89:ARG:HD2	1:B:95:GLU:OE1	2.13	0.49
1:B:213:HIS:HB2	1:B:236:ARG:NH2	2.27	0.49
1:B:86:THR:HG22	1:B:91:LEU:HG	1.95	0.49
1:A:259:TRP:NE1	4:A:580:HOH:O	2.43	0.49
1:A:176:ARG:CB	1:A:182:MET:HE3	2.42	0.49
1:A:217:LEU:HD12	1:A:232:LEU:CG	2.40	0.49
1:A:79:VAL:HG11	1:A:365:ALA:HB3	1.94	0.49
1:A:159:LEU:HD13	1:A:161:PHE:CZ	2.48	0.49
1:A:314:VAL:CG1	1:A:318:ARG:NH1	2.74	0.49
1:B:164:HIS:O	1:B:165:PRO:C	2.56	0.49
1:A:51:ARG:NH2	1:A:199:PRO:O	2.46	0.48
1:B:228:ASP:HB3	1:B:231:MET:HB3	1.95	0.48
1:B:308:PHE:O	1:B:309:ALA:C	2.56	0.48
1:A:162:ASP:O	1:A:163:ARG:HB2	2.13	0.48
1:A:222:ARG:H	1:A:222:ARG:HD3	1.79	0.48
1:B:118:ALA:O	1:B:122:THR:HG22	2.12	0.48
1:A:62:LYS:HE2	1:A:163:ARG:NH1	2.28	0.48
1:A:123:SER:HB3	1:B:122:THR:CG2	2.44	0.48
1:A:219:LYS:O	1:A:220:ARG:HB3	2.13	0.48
1:A:183:THR:HG22	1:A:185:GLN:N	2.28	0.48
1:B:107:LEU:CA	1:B:112:ILE:HG22	2.43	0.48
1:A:295:PRO:HA	1:A:298:LEU:HD12	1.95	0.47
1:B:187:VAL:O	1:B:191:VAL:HG23	2.15	0.47
1:A:177:TYR:HA	1:A:182:MET:O	2.14	0.47
1:A:332:SER:HB2	1:A:333:PRO:HD2	1.95	0.47
1:B:60:MET:HE1	1:B:206:GLY:HA3	1.96	0.47
1:A:79:VAL:HG11	1:A:365:ALA:CB	2.45	0.47
1:B:78:ILE:HA	1:B:158:THR:O	2.14	0.47
1:A:219:LYS:NZ	1:A:220:ARG:O	2.41	0.47
1:A:310:TRP:O	1:A:313:ASP:HB2	2.13	0.47
1:B:166:ILE:HG21	1:B:322:MET:HE1	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:PRO:O	1:A:335:GLY:N	2.41	0.47
1:A:221:GLN:NE2	1:A:222:ARG:NH1	2.62	0.47
1:B:125:GLN:HE22	3:B:2:PE2:C6	2.28	0.47
1:A:294:ALA:HA	1:A:295:PRO:HD2	1.77	0.47
3:B:2:PE2:H3'1	3:B:2:PE2:H1'1	1.61	0.46
1:B:208:LEU:O	1:B:236:ARG:NH2	2.42	0.46
1:A:153:PRO:HB3	1:A:349:GLN:HE22	1.81	0.46
1:B:241:LEU:HD13	1:B:288:LEU:HA	1.98	0.46
1:B:319:LEU:HD23	1:B:322:MET:SD	2.55	0.46
1:B:240:GLY:O	1:B:243:ALA:N	2.48	0.46
1:B:262:LEU:HD13	1:B:284:ILE:HG23	1.98	0.46
1:A:164:HIS:ND1	1:A:165:PRO:HD2	2.30	0.46
1:B:62:LYS:CE	1:B:163:ARG:NH1	2.76	0.46
1:B:107:LEU:HD23	1:B:112:ILE:HG23	1.97	0.46
1:A:142:HIS:HE1	4:A:515:HOH:O	1.98	0.46
1:B:362:CYS:SG	1:B:366:ARG:CZ	3.04	0.46
1:A:126:ILE:CD1	1:B:126:ILE:HD11	2.45	0.45
1:A:221:GLN:NE2	1:A:222:ARG:HH12	2.14	0.45
1:A:48:THR:OG1	1:A:155:PRO:HA	2.16	0.45
1:A:121:MET:O	1:A:125:GLN:HG2	2.15	0.45
1:B:248:TYR:CD2	1:B:249:LEU:HD23	2.52	0.45
1:A:50:LEU:HD23	1:A:349:GLN:HG2	1.98	0.45
1:A:221:GLN:HB2	4:A:627:HOH:O	2.16	0.45
1:B:50:LEU:HD12	1:B:158:THR:HG23	1.97	0.45
1:B:56:GLY:O	1:B:163:ARG:NH2	2.49	0.45
1:B:164:HIS:CD2	1:B:203:ILE:HD11	2.51	0.45
1:B:240:GLY:O	1:B:241:LEU:C	2.59	0.45
1:A:79:VAL:HB	1:A:159:LEU:HD23	1.99	0.45
1:B:359:PRO:O	1:B:362:CYS:HB3	2.17	0.45
1:B:121:MET:HB3	1:B:182:MET:CE	2.45	0.45
1:A:183:THR:O	1:A:187:VAL:HG23	2.17	0.45
1:B:297:LEU:HD22	1:B:308:PHE:CD1	2.52	0.45
1:A:109:GLN:O	1:A:111:GLU:N	2.50	0.45
1:B:58:HIS:HB3	3:B:2:PE2:H3'1	1.99	0.45
1:B:159:LEU:HD11	1:B:361:ILE:CG2	2.47	0.45
1:A:121:MET:HE3	1:A:176:ARG:HD2	1.98	0.45
1:A:339:ALA:O	1:A:343:LEU:HG	2.17	0.45
1:B:219:LYS:O	1:B:220:ARG:CB	2.65	0.45
1:B:235:ILE:HD12	1:B:235:ILE:HA	1.81	0.44
1:A:283:HIS:HB3	1:A:286:ASP:OD2	2.17	0.44
1:B:102:THR:O	1:B:105:HIS:HB3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:TRP:O	1:B:262:LEU:HB2	2.18	0.44
1:A:78:ILE:CG1	1:A:158:THR:HB	2.42	0.44
1:B:186:ALA:O	1:B:189:ALA:HB3	2.17	0.44
1:B:80:TYR:O	1:B:82:PRO:HD3	2.18	0.44
1:B:67:GLN:O	1:B:70:VAL:N	2.50	0.44
1:B:176:ARG:CB	1:B:182:MET:HE3	2.46	0.44
1:B:176:ARG:HB3	1:B:182:MET:HE3	1.99	0.44
1:B:248:TYR:CZ	1:B:253:GLY:HA3	2.53	0.44
1:A:48:THR:HA	1:A:350:THR:O	2.17	0.43
1:A:372:MET:HE1	1:B:307:VAL:HG13	2.00	0.43
1:B:148:GLY:O	1:B:152:ALA:N	2.50	0.43
1:B:205:LEU:HB2	1:B:326:ILE:HD13	1.99	0.43
1:B:232:LEU:HD23	1:B:232:LEU:HA	1.82	0.43
1:A:210:GLU:O	1:A:214:ILE:HG13	2.18	0.43
1:B:101:TYR:CD2	1:B:225:GLU:HA	2.53	0.43
1:B:140:ALA:HB3	1:B:141:PRO:HD3	2.00	0.43
1:B:208:LEU:HD12	1:B:209:PRO:CD	2.47	0.43
1:A:95:GLU:O	1:A:96:THR:C	2.60	0.43
1:A:217:LEU:HD12	1:A:232:LEU:CD2	2.49	0.43
1:B:146:GLU:HG2	1:B:350:THR:HG22	2.01	0.43
1:A:368:PHE:CD1	1:A:368:PHE:C	2.96	0.43
1:B:124:ALA:O	1:B:128:MET:HG2	2.19	0.43
1:B:67:GLN:O	1:B:68:LEU:C	2.61	0.43
1:B:234:ALA:O	1:B:237:ARG:HB3	2.19	0.43
1:A:221:GLN:HE22	1:A:222:ARG:HH12	1.66	0.42
1:B:101:TYR:CZ	1:B:225:GLU:HG3	2.54	0.42
1:B:210:GLU:HG3	1:B:214:ILE:CD1	2.46	0.42
1:B:288:LEU:O	1:B:289:PHE:C	2.62	0.42
1:A:219:LYS:HG2	1:A:220:ARG:H	1.84	0.42
1:B:216:ARG:O	1:B:217:LEU:C	2.60	0.42
1:A:126:ILE:HD13	1:A:190:PHE:CE1	2.52	0.42
1:A:130:MET:HA	4:B:630:HOH:O	2.19	0.42
1:A:257:GLU:HG3	1:A:258:ASP:OD1	2.19	0.42
1:B:77:ASP:O	1:B:157:LEU:HD12	2.18	0.42
1:A:188:LEU:HD23	1:A:188:LEU:HA	1.90	0.42
1:A:314:VAL:O	1:A:315:LEU:C	2.63	0.42
1:B:295:PRO:HG2	1:B:296:GLU:OE1	2.20	0.42
1:A:62:LYS:HA	1:A:329:TYR:OH	2.20	0.42
1:B:153:PRO:HA	1:B:154:PRO:HA	1.80	0.42
1:B:213:HIS:CB	1:B:236:ARG:HH21	2.33	0.42
1:A:62:LYS:O	1:A:66:THR:HG22	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:TYR:HB2	1:B:308:PHE:CD2	2.55	0.42
1:A:79:VAL:CG1	1:A:80:TYR:H	2.33	0.42
1:A:89:ARG:HD2	1:A:95:GLU:CD	2.45	0.42
1:A:200:GLY:HA3	1:A:350:THR:CG2	2.50	0.42
1:A:299:ALA:O	1:A:302:GLY:N	2.47	0.42
1:B:51:ARG:NH1	1:B:53:TYR:OH	2.53	0.42
1:B:89:ARG:O	1:B:94:SER:HA	2.20	0.42
1:B:213:HIS:CD2	1:B:217:LEU:CD2	3.01	0.42
1:A:79:VAL:CG1	1:A:80:TYR:N	2.82	0.42
1:A:311:ALA:O	1:A:314:VAL:HB	2.20	0.42
1:A:133:ALA:O	1:A:134:VAL:C	2.63	0.41
1:A:262:LEU:HD21	1:A:284:ILE:HG23	2.02	0.41
1:B:296:GLU:OE1	1:B:296:GLU:N	2.52	0.41
1:A:368:PHE:CE2	1:A:372:MET:HE3	2.55	0.41
1:A:89:ARG:HA	1:A:95:GLU:HA	2.01	0.41
1:A:200:GLY:HA3	1:A:350:THR:HG23	2.02	0.41
1:B:193:LEU:HD23	1:B:193:LEU:HA	1.58	0.41
1:A:213:HIS:CE1	1:A:232:LEU:CD2	3.02	0.41
1:A:123:SER:HB3	1:B:122:THR:HG23	2.03	0.41
1:B:65:THR:HG22	1:B:336:CYS:HB2	2.02	0.41
1:B:182:MET:HE3	1:B:182:MET:HB2	1.61	0.41
1:A:217:LEU:O	1:A:221:GLN:HG3	2.20	0.41
1:A:262:LEU:O	1:A:293:ARG:NH1	2.54	0.41
1:B:52:VAL:HG22	1:B:347:MET:SD	2.61	0.41
1:B:154:PRO:HB2	1:B:155:PRO:CD	2.51	0.41
1:B:217:LEU:HD13	1:B:217:LEU:HA	1.75	0.41
1:B:112:ILE:HG12	1:B:116:ASP:CB	2.51	0.41
1:A:142:HIS:CE1	4:A:515:HOH:O	2.74	0.40
1:A:255:TRP:O	1:A:259:TRP:HB2	2.21	0.40
1:A:325:PHE:CD1	1:A:325:PHE:N	2.89	0.40
1:B:255:TRP:HB2	1:B:284:ILE:HD11	2.01	0.40
1:B:203:ILE:O	1:B:324:VAL:HA	2.21	0.40
1:B:69:LEU:HD22	1:B:340:LEU:HD21	2.04	0.40
1:B:85:MET:CE	1:B:88:TRP:HB2	2.52	0.40
1:B:122:THR:HG23	1:B:123:SER:N	2.37	0.40
1:B:297:LEU:HD22	1:B:308:PHE:CE1	2.56	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	301/331 (91%)	260 (86%)	31 (10%)	10 (3%)	3	2
1	B	303/331 (92%)	267 (88%)	31 (10%)	5 (2%)	7	9
All	All	604/662 (91%)	527 (87%)	62 (10%)	15 (2%)	4	4

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	222	ARG
1	B	149	SER
1	B	220	ARG
1	A	110	GLY
1	A	220	ARG
1	A	259	TRP
1	A	344	THR
1	A	163	ARG
1	A	333	PRO
1	B	94	SER
1	B	221	GLN
1	A	140	ALA
1	B	154	PRO
1	A	153	PRO
1	A	82	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	245/264 (93%)	218 (89%)	27 (11%)	6	8
1	B	246/264 (93%)	207 (84%)	39 (16%)	2	3
All	All	491/528 (93%)	425 (87%)	66 (13%)	4	4

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	ASP
1	A	62	LYS
1	A	64	THR
1	A	78	ILE
1	A	81	VAL
1	A	107	LEU
1	A	119	VAL
1	A	130	MET
1	A	134	VAL
1	A	165	PRO
1	A	178	LEU
1	A	196	PRO
1	A	197	THR
1	A	203	ILE
1	A	217	LEU
1	A	227	LEU
1	A	232	LEU
1	A	293	ARG
1	A	300	PRO
1	A	307	VAL
1	A	320	ARG
1	A	337	ARG
1	A	338	ASP
1	A	348	VAL
1	A	362	CYS
1	A	368	PHE
1	A	372	MET
1	B	49	LEU
1	B	50	LEU
1	B	62	LYS
1	B	72	LEU
1	B	85	MET
1	B	107	LEU
1	B	109	GLN
1	B	112	ILE

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Mol	Chain	Res	Type
1	B	119	VAL
1	B	130	MET
1	B	134	VAL
1	B	138	VAL
1	B	154	PRO
1	B	165	PRO
1	B	173	PRO
1	B	178	LEU
1	B	188	LEU
1	B	209	PRO
1	B	212	ARG
1	B	217	LEU
1	B	220	ARG
1	B	221	GLN
1	B	222	ARG
1	B	226	ARG
1	B	227	LEU
1	B	232	LEU
1	B	247	ARG
1	B	249	LEU
1	B	262	LEU
1	B	307	VAL
1	B	318	ARG
1	B	333	PRO
1	B	341	LEU
1	B	349	GLN
1	B	352	VAL
1	B	354	THR
1	B	360	THR
1	B	372	MET
1	B	374	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	HIS
1	A	67	GLN
1	A	142	HIS
1	A	221	GLN
1	A	349	GLN
1	A	351	HIS
1	B	142	HIS

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Mol	Chain	Res	Type
1	B	301	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](#)

4 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	PE2	B	2	-	19,19,19	0.77	0	26,26,26	2.00	8 (30%)
2	SO4	B	4	-	4,4,4	0.67	0	6,6,6	1.27	0
2	SO4	A	3	-	4,4,4	0.57	0	6,6,6	1.17	1 (16%)
3	PE2	A	1	-	19,19,19	1.17	1 (5%)	26,26,26	2.98	7 (26%)

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	PE2	A	1	-	-	6/9/9/9	0/2/2/2
3	PE2	B	2	-	-	1/9/9/9	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1	PE2	C4-N3	2.31	1.39	1.34

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1	PE2	C12-C1'-N9	11.25	130.78	112.41
3	A	1	PE2	C1'-C12-C2'	-6.27	100.99	114.85
3	B	2	PE2	C12-C1'-N9	-5.13	104.03	112.41
3	A	1	PE2	N9-C8-N7	4.92	118.65	113.52
3	B	2	PE2	N9-C8-N7	3.28	116.94	113.52
3	B	2	PE2	N2-C2-N3	-3.02	113.77	119.67
3	B	2	PE2	C5-C6-N1	-2.93	105.80	113.25
3	A	1	PE2	C8-N9-C4	-2.61	103.07	105.89
3	A	1	PE2	O6-C6-N1	2.61	125.02	120.11
3	A	1	PE2	O4'-C4'-C2'	-2.51	99.67	112.36
3	B	2	PE2	N9-C4-N3	-2.44	121.89	125.46
2	A	3	SO4	O4-S-O2	2.35	121.83	109.56
3	B	2	PE2	C2-N3-C4	-2.34	108.27	112.30
3	B	2	PE2	C8-N9-C4	-2.32	103.39	105.89
3	A	1	PE2	C1'-N9-C4	2.31	130.89	127.13
3	B	2	PE2	O6-C6-C5	2.17	132.28	126.53

There are no chirality outliers.

All (7) torsion outliers are listed below:

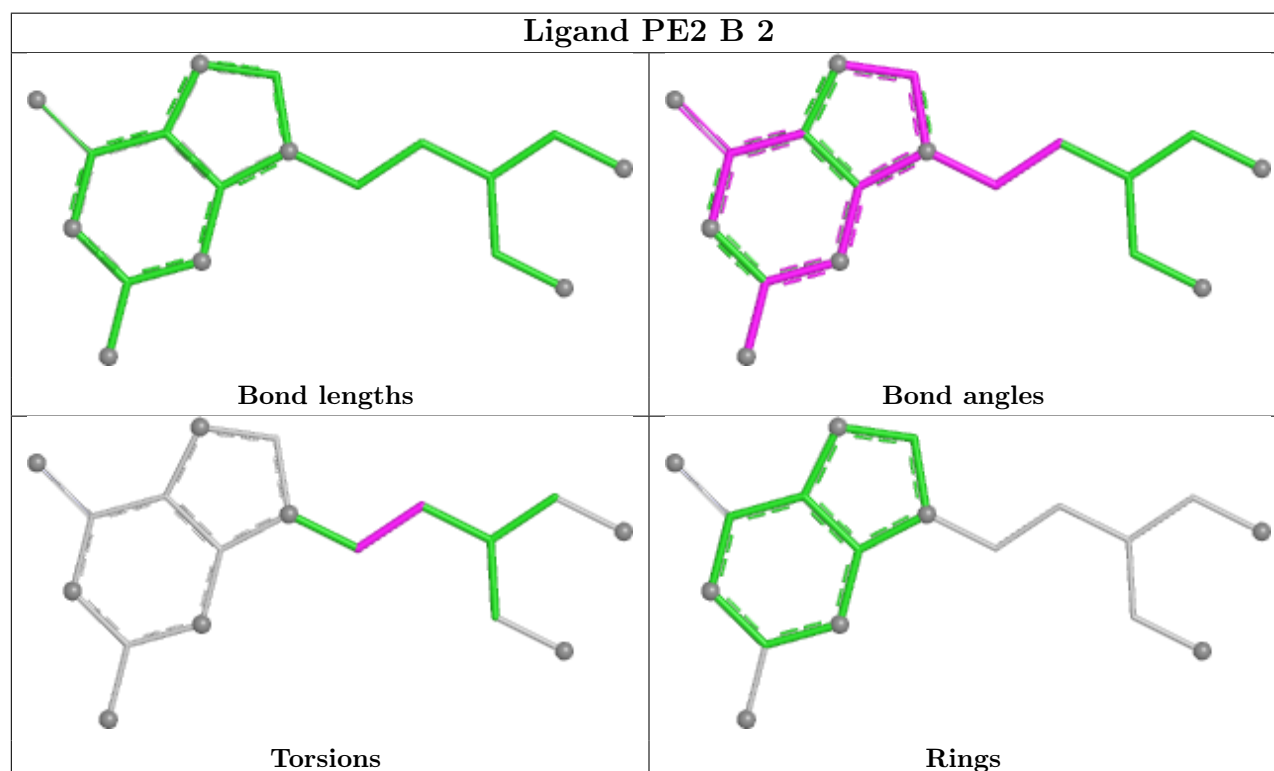
Mol	Chain	Res	Type	Atoms
3	A	1	PE2	C12-C1'-N9-C8
3	A	1	PE2	C12-C1'-N9-C4
3	B	2	PE2	N9-C1'-C12-C2'
3	A	1	PE2	C4'-C2'-C3'-O3'
3	A	1	PE2	C1'-C12-C2'-C4'
3	A	1	PE2	C1'-C12-C2'-C3'
3	A	1	PE2	C12-C2'-C3'-O3'

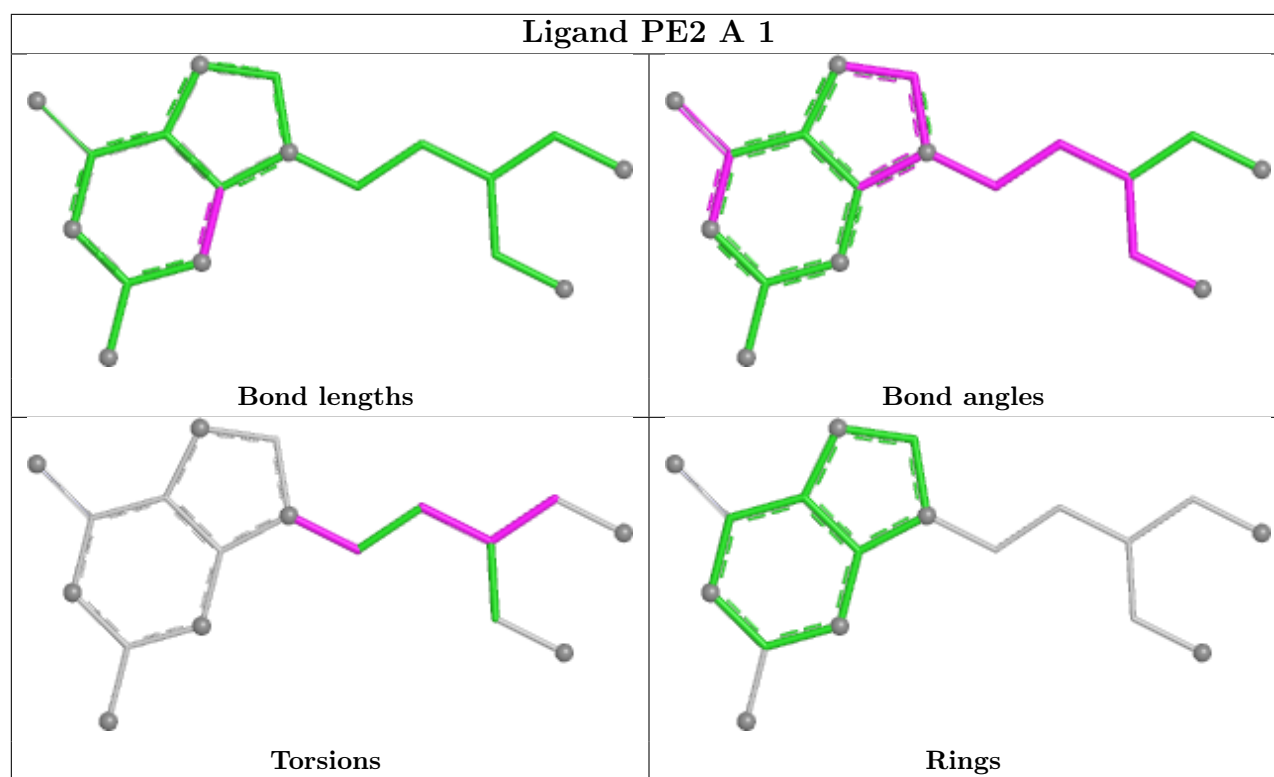
There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2	PE2	5	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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	309/331 (93%)	-1.59	0 100 100	5, 20, 42, 64	0
1	B	311/331 (93%)	-1.58	0 100 100	5, 19, 39, 52	0
All	All	620/662 (93%)	-1.58	0 100 100	5, 19, 40, 64	0

There are no RSRZ outliers to report.

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.

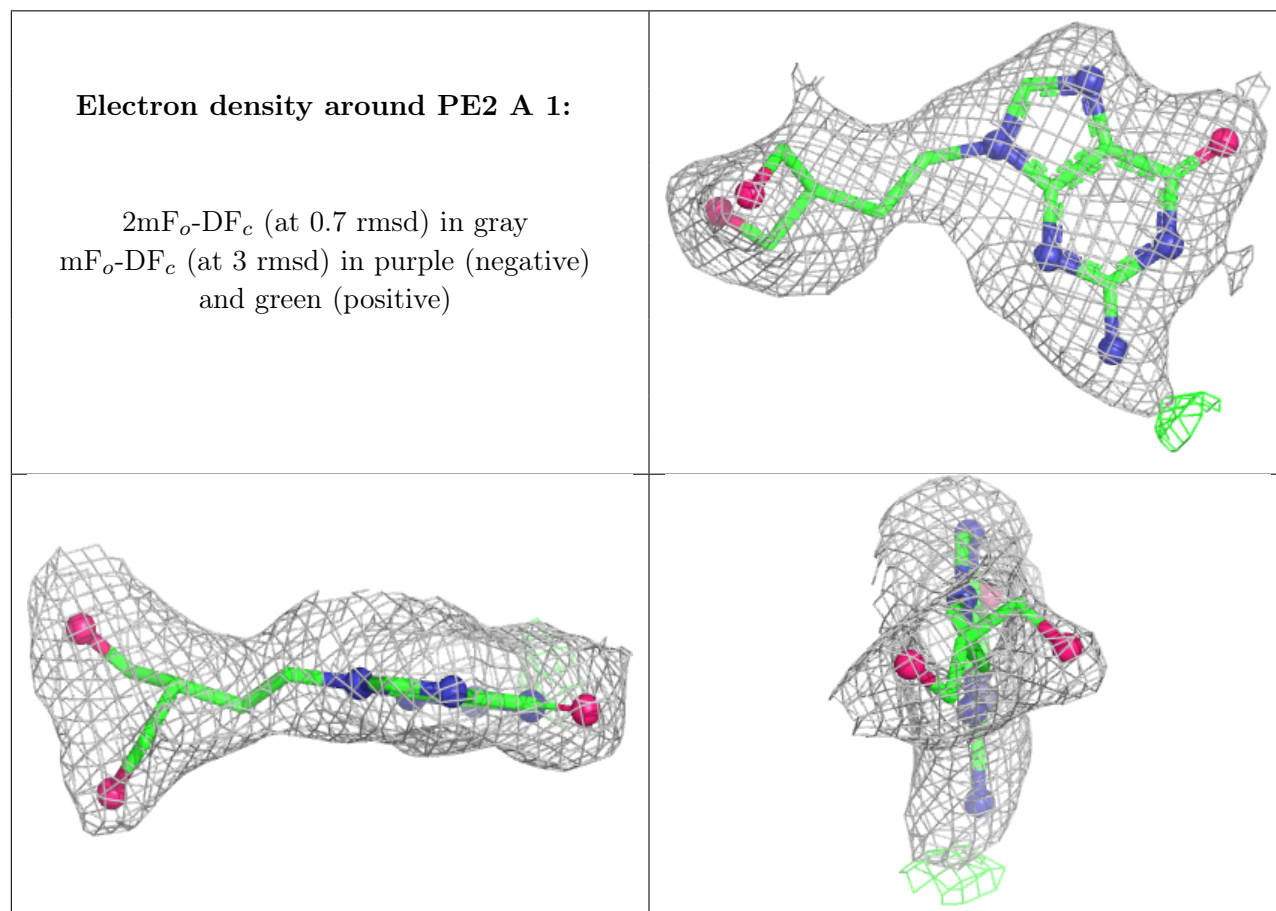
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
2	SO4	A	3	5/5	1.00	0.02	23,24,29,30	0
2	SO4	B	4	5/5	1.00	0.02	17,18,19,21	0
3	PE2	A	1	18/18	1.00	0.02	18,21,23,23	0
3	PE2	B	2	18/18	1.00	0.02	18,23,38,39	0

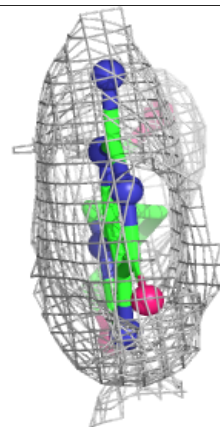
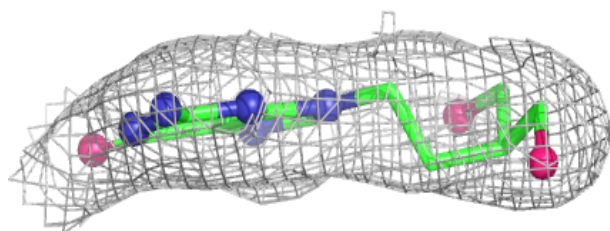
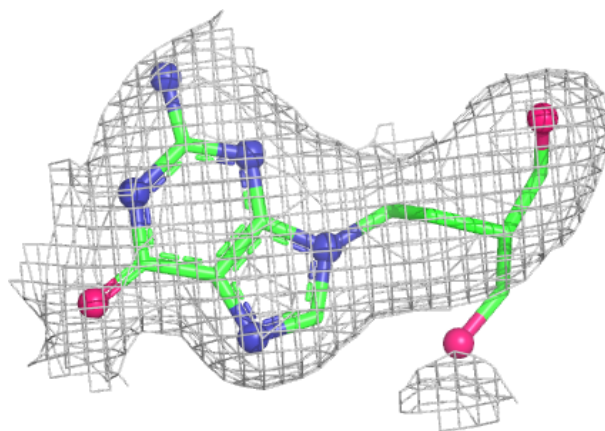
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

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 PE2 B 2:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.