



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:00 PM UTC

PDB ID : 1ODJ / pdb_00001odj
Title : PURINE NUCLEOSIDE PHOSPHORYLASE FROM THERMUS THERMOPHILUS
Authors : Tahirov, T.H.; Inagaki, E.; Miyano, M.
Deposited on : 2003-02-19
Resolution : 2.40 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

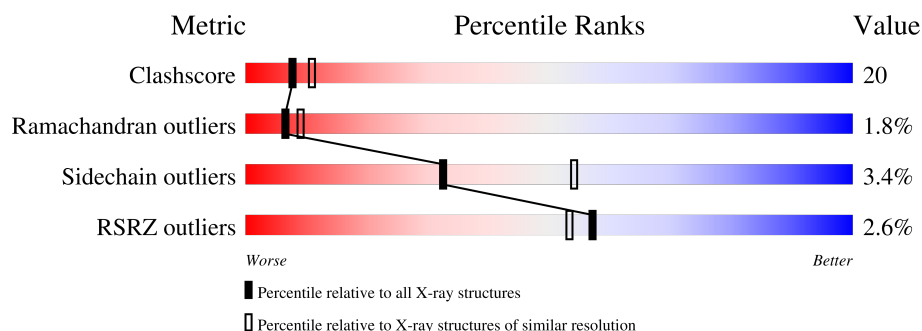
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	235	3% 64% 32% .
1	B	235	% 62% 33% 5%
1	C	235	3% 53% 42% .
1	D	235	3% 56% 37% 7%
1	E	235	3% 66% 31% .
1	F	235	2% 72% 25% .

2 Entry composition [i](#)

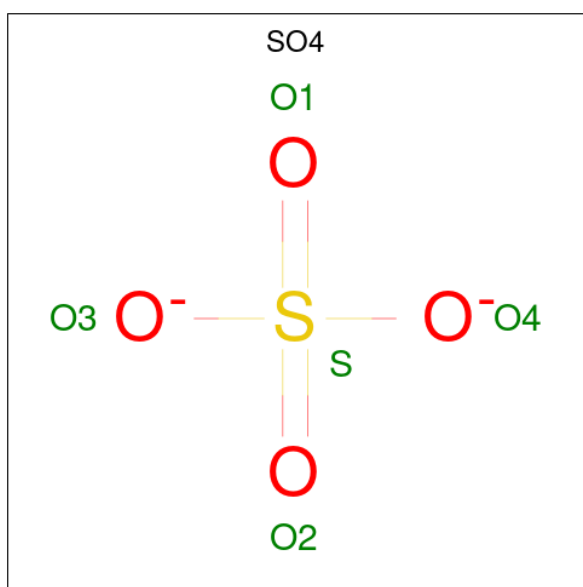
There are 4 unique types of molecules in this entry. The entry contains 11124 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PURINE NUCLEOSIDE PHOSPHORYLASE.

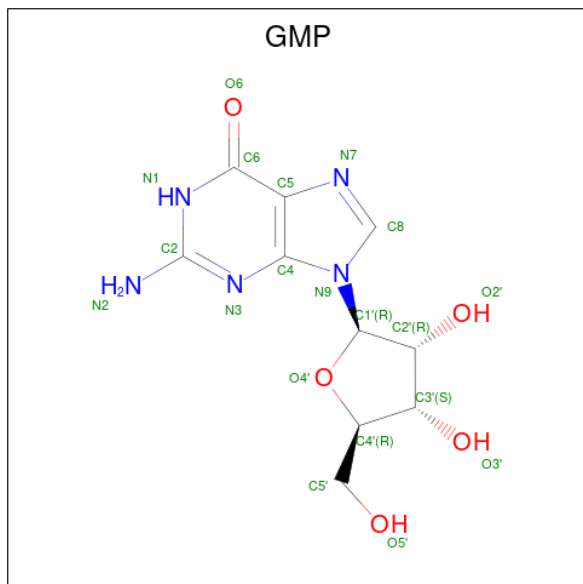
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	0	0	0
			1786	1133	325	324	4			
1	B	234	Total	C	N	O	S	0	0	0
			1786	1133	325	324	4			
1	C	234	Total	C	N	O	S	0	0	0
			1786	1133	325	324	4			
1	D	234	Total	C	N	O	S	0	0	0
			1786	1133	325	324	4			
1	E	234	Total	C	N	O	S	0	0	0
			1786	1133	325	324	4			
1	F	234	Total	C	N	O	S	0	0	0
			1786	1133	325	324	4			

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



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

- Molecule 3 is GUANOSINE (CCD ID: GMP) (formula: $C_{10}H_{13}N_5O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 20 10 5 5	0	0
3	B	1	Total C N O 20 10 5 5	0	0
3	C	1	Total C N O 20 10 5 5	0	0
3	D	1	Total C N O 20 10 5 5	0	0
3	E	1	Total C N O 20 10 5 5	0	0
3	F	1	Total C N O 20 10 5 5	0	0

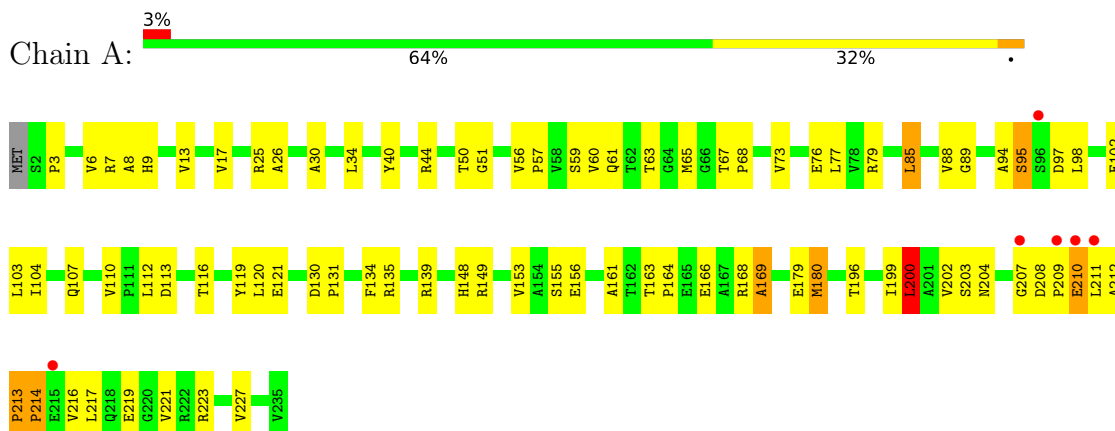
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	58	Total 58	O 58	0	0
4	B	50	Total 50	O 50	0	0
4	C	27	Total 27	O 27	0	0
4	D	23	Total 23	O 23	0	0
4	E	40	Total 40	O 40	0	0
4	F	60	Total 60	O 60	0	0

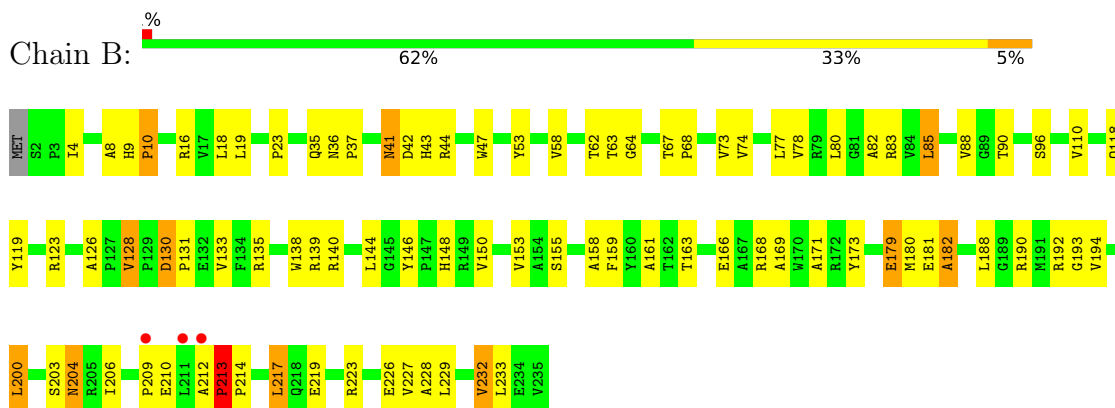
3 Residue-property plots

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.

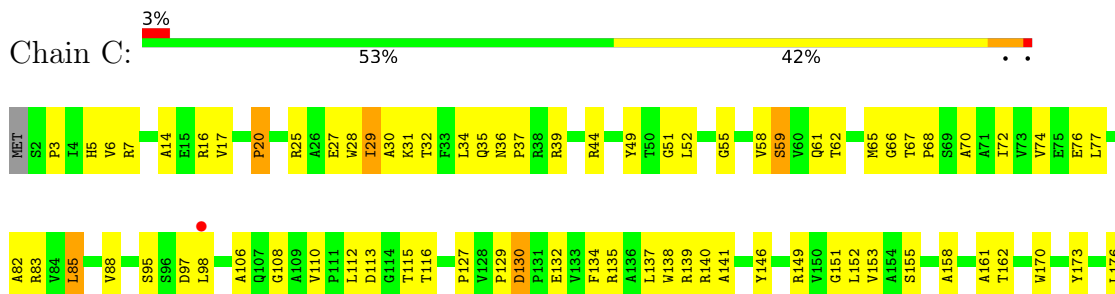
• Molecule 1: PURINE NUCLEOSIDE PHOSPHORYLASE

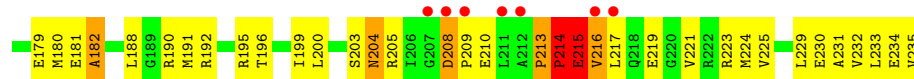


• Molecule 1: PURINE NUCLEOSIDE PHOSPHORYLASE

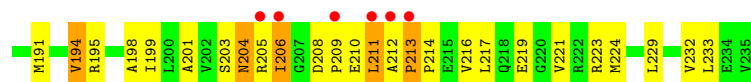


• Molecule 1: PURINE NUCLEOSIDE PHOSPHORYLASE

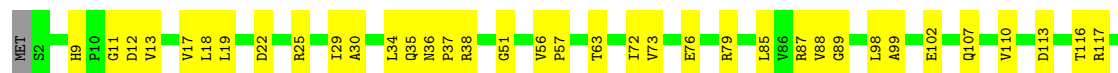




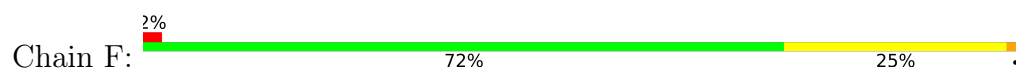
● Molecule 1: PURINE NUCLEOSIDE PHOSPHORYLASE



● Molecule 1: PURINE NUCLEOSIDE PHOSPHORYLASE



● Molecule 1: PURINE NUCLEOSIDE PHOSPHORYLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	132.76Å 132.76Å 172.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.93 – 2.40 14.93 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.8 (14.93-2.40) 96.4 (14.93-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.37 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.286 0.228 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.189	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 64.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	11124	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.12% 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

5.1 Standard geometry

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

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	0.53	0/1828	1.04	14/2493 (0.6%)
1	B	0.47	0/1828	1.01	9/2493 (0.4%)
1	C	0.46	0/1828	0.96	10/2493 (0.4%)
1	D	0.45	0/1828	1.01	14/2493 (0.6%)
1	E	0.50	0/1828	0.98	3/2493 (0.1%)
1	F	0.53	0/1828	1.04	8/2493 (0.3%)
All	All	0.49	0/10968	1.01	58/14958 (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	1

There are no bond length outliers.

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	213	PRO	CA-C-N	7.94	129.77	119.84
1	A	213	PRO	C-N-CA	7.94	129.77	119.84
1	B	153	VAL	N-CA-C	-7.89	97.76	108.96
1	D	130	ASP	N-CA-C	-7.10	101.12	109.93
1	F	56	VAL	N-CA-C	6.96	114.62	108.63
1	D	32	THR	N-CA-C	6.96	118.65	111.14
1	B	180	MET	N-CA-C	6.76	119.53	110.88
1	D	123	ARG	N-CA-C	6.62	118.24	109.83
1	D	64	GLY	N-CA-C	-6.61	102.92	112.17
1	A	130	ASP	N-CA-C	-6.49	101.48	109.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	110	VAL	CA-C-N	6.47	127.92	119.84
1	B	110	VAL	C-N-CA	6.47	127.92	119.84
1	A	180	MET	N-CA-C	6.39	118.60	110.61
1	F	19	LEU	N-CA-C	6.33	120.86	108.21
1	B	4	ILE	N-CA-C	6.29	116.46	110.42
1	B	148	HIS	N-CA-C	6.29	116.58	108.34
1	C	130	ASP	N-CA-C	-6.29	102.13	109.93
1	D	182	ALA	N-CA-C	6.11	118.44	111.11
1	C	219	GLU	N-CA-C	-6.00	104.66	111.14
1	F	180	MET	N-CA-C	5.83	119.61	111.56
1	A	153	VAL	N-CA-C	-5.82	100.70	108.96
1	C	208	ASP	CA-C-N	5.80	127.09	119.84
1	C	208	ASP	C-N-CA	5.80	127.09	119.84
1	A	169	ALA	N-CA-C	-5.78	104.89	111.07
1	E	17	VAL	N-CA-C	5.75	116.57	108.17
1	A	26	ALA	N-CA-C	-5.74	104.71	110.97
1	A	148	HIS	N-CA-C	5.62	115.97	108.38
1	A	208	ASP	CA-C-N	5.54	125.21	119.56
1	A	208	ASP	C-N-CA	5.54	125.21	119.56
1	D	208	ASP	CA-C-N	5.50	125.59	119.87
1	D	208	ASP	C-N-CA	5.50	125.59	119.87
1	B	41	ASN	N-CA-C	5.49	117.44	108.55
1	F	230	GLU	N-CA-C	-5.44	105.25	111.07
1	D	126	ALA	CA-C-N	5.44	126.64	119.84
1	D	126	ALA	C-N-CA	5.44	126.64	119.84
1	F	112	LEU	N-CA-C	-5.35	106.62	112.72
1	C	210	GLU	N-CA-C	5.34	116.11	108.74
1	C	215	GLU	N-CA-C	5.32	117.08	111.28
1	A	200	LEU	N-CA-C	5.31	117.45	109.23
1	D	183	SER	N-CA-C	5.29	116.73	111.07
1	E	173	TYR	N-CA-C	-5.26	106.09	112.88
1	A	79	ARG	N-CA-C	-5.25	105.67	111.71
1	C	141	ALA	N-CA-C	-5.23	105.22	111.03
1	D	180	MET	N-CA-C	5.22	118.70	111.71
1	C	29	ILE	N-CA-C	-5.20	105.78	110.82
1	B	64	GLY	N-CA-C	-5.16	104.94	112.17
1	D	107	GLN	N-CA-C	-5.14	106.51	112.89
1	A	9	HIS	CA-C-N	5.14	126.26	119.84
1	A	9	HIS	C-N-CA	5.14	126.26	119.84
1	F	152	LEU	N-CA-C	5.12	118.10	110.52
1	C	182	ALA	N-CA-C	5.12	117.25	111.11
1	B	169	ALA	N-CA-C	-5.12	105.61	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(^o)	Ideal(^o)
1	E	180	MET	N-CA-C	5.12	118.56	111.71
1	F	210	GLU	N-CA-C	5.08	114.33	108.49
1	C	106	ALA	N-CA-C	5.08	116.46	109.15
1	F	200	LEU	N-CA-C	5.04	117.79	109.72
1	D	110	VAL	CA-C-N	5.01	126.11	119.84
1	D	110	VAL	C-N-CA	5.01	126.11	119.84

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	119	TYR	Sidechain

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	1786	0	1792	57	0
1	B	1786	0	1792	72	0
1	C	1786	0	1792	104	0
1	D	1786	0	1792	97	0
1	E	1786	0	1792	71	0
1	F	1786	0	1792	49	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	1	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	20	0	13	1	0
3	B	20	0	13	2	0
3	C	20	0	13	2	0
3	D	20	0	13	4	0
3	E	20	0	13	3	0
3	F	20	0	13	1	0
4	A	58	0	0	4	0
4	B	50	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	27	0	0	4	0
4	D	23	0	0	0	0
4	E	40	0	0	4	0
4	F	60	0	0	5	0
All	All	11124	0	10830	433	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 20.

All (433) 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:212:ALA:HB1	1:A:213:PRO:HD2	1.32	1.09
1:C:20:PRO:HD2	1:C:61:GLN:O	1.70	0.91
1:F:212:ALA:HA	4:F:2058:HOH:O	1.68	0.91
1:F:3:PRO:HB3	1:F:76:GLU:HG2	1.50	0.91
1:E:88:VAL:HB	1:E:224:MET:HE3	1.51	0.90
1:B:203:SER:HB2	1:B:217:LEU:HD11	1.52	0.88
1:C:205:ARG:HB2	4:C:2023:HOH:O	1.79	0.81
1:F:213:PRO:HB2	1:F:216:VAL:HG23	1.62	0.81
1:C:191:MET:HE2	1:C:192:ARG:HH22	1.44	0.81
1:A:95:SER:HB2	1:A:97:ASP:OD1	1.82	0.79
1:D:88:VAL:HB	1:D:224:MET:HE3	1.65	0.79
1:E:37:PRO:O	1:E:38:ARG:HD3	1.84	0.78
1:E:212:ALA:HB1	1:E:213:PRO:CD	2.14	0.78
1:D:36:ASN:N	1:D:36:ASN:HD22	1.82	0.78
1:C:44:ARG:HG2	1:D:22:ASP:HB2	1.65	0.77
1:E:147:PRO:HD2	1:E:223:ARG:NH1	1.99	0.77
1:B:212:ALA:HB1	1:B:213:PRO:HD2	1.67	0.76
1:D:132:GLU:HG3	1:D:195:ARG:HH12	1.49	0.76
1:C:16:ARG:HD3	1:C:232:VAL:HB	1.68	0.76
1:D:178:PHE:CD1	1:D:206:ILE:HD11	2.20	0.76
1:C:85:LEU:N	1:C:85:LEU:HD12	2.01	0.75
1:F:98:LEU:HD21	1:F:149:ARG:HG3	1.68	0.75
1:C:25:ARG:O	1:C:29:ILE:HG13	1.86	0.75
1:E:212:ALA:HB1	1:E:213:PRO:HD2	1.69	0.74
1:A:98:LEU:HD21	1:A:104:ILE:HD11	1.70	0.73
1:D:209:PRO:O	1:D:210:GLU:HG3	1.88	0.73
1:D:212:ALA:HB1	1:D:213:PRO:HD2	1.70	0.72
1:F:212:ALA:HB1	1:F:213:PRO:CD	2.19	0.72
1:F:130:ASP:OD2	1:F:133:VAL:HG23	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:ARG:HH12	1:A:89:GLY:HA2	1.54	0.71
1:B:123:ARG:HD3	4:B:2029:HOH:O	1.91	0.71
1:C:34:LEU:HB3	1:C:51:GLY:HA3	1.72	0.71
1:B:9:HIS:HB3	1:B:10:PRO:HD2	1.73	0.71
1:D:78:VAL:HG22	1:D:194:VAL:HG11	1.72	0.70
1:A:210:GLU:O	1:A:211:LEU:HD23	1.92	0.69
1:C:82:ALA:O	1:C:83:ARG:NH1	2.23	0.69
1:E:204:ASN:HD22	1:E:205:ARG:N	1.90	0.69
1:D:232:VAL:HG23	1:D:233:LEU:HD12	1.74	0.69
1:C:113:ASP:OD1	1:C:116:THR:HG23	1.93	0.68
1:A:85:LEU:N	1:A:85:LEU:HD12	2.08	0.68
1:E:178:PHE:CD1	1:E:206:ILE:HD11	2.29	0.67
1:D:36:ASN:N	1:D:37:PRO:HD3	2.10	0.67
1:D:85:LEU:HD12	1:D:85:LEU:N	2.10	0.67
1:C:221:VAL:O	1:C:225:VAL:HG23	1.94	0.67
1:E:147:PRO:HD2	1:E:223:ARG:HH12	1.58	0.67
1:A:212:ALA:HB1	1:A:213:PRO:CD	2.16	0.66
1:D:212:ALA:HB1	1:D:213:PRO:CD	2.25	0.66
1:A:98:LEU:HD11	1:A:149:ARG:HG3	1.77	0.65
1:D:139:ARG:HE	1:E:139:ARG:HH22	1.43	0.65
1:F:3:PRO:HB2	1:F:6:VAL:O	1.97	0.65
1:F:25:ARG:O	1:F:29:ILE:HG13	1.96	0.65
1:F:213:PRO:HB2	1:F:216:VAL:CG2	2.27	0.64
1:C:61:GLN:NE2	1:C:62:THR:H	1.95	0.64
1:B:190:ARG:HD3	4:B:2042:HOH:O	1.98	0.64
1:D:179:GLU:CD	1:D:182:ALA:H	2.06	0.64
1:D:217:LEU:O	1:D:221:VAL:HG23	1.97	0.64
1:D:211:LEU:HD13	1:D:217:LEU:CD1	2.28	0.63
1:D:93:ALA:O	1:D:206:ILE:HG13	1.98	0.63
1:C:203:SER:O	1:C:204:ASN:HB3	1.97	0.63
1:E:146:TYR:HB3	4:E:2037:HOH:O	1.98	0.63
1:B:163:THR:HG23	1:B:166:GLU:OE1	1.99	0.63
1:C:98:LEU:CD2	1:C:149:ARG:HG3	2.28	0.63
1:D:132:GLU:HG3	1:D:195:ARG:NH1	2.14	0.62
1:A:25:ARG:NH1	1:A:89:GLY:HA2	2.14	0.62
1:B:126:ALA:O	1:B:128:VAL:HG22	1.99	0.62
1:B:163:THR:OG1	1:B:166:GLU:HG3	1.99	0.62
1:B:212:ALA:HB1	1:B:213:PRO:CD	2.29	0.62
1:C:179:GLU:OE1	1:C:182:ALA:HB3	2.00	0.62
1:D:204:ASN:C	1:D:204:ASN:HD22	2.09	0.61
1:D:128:VAL:HB	1:E:107:GLN:NE2	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:VAL:HG22	1:A:59:SER:HB2	1.83	0.61
3:C:1237:GMP:N3	3:C:1237:GMP:H2'	2.15	0.61
1:B:140:ARG:HD3	4:B:2034:HOH:O	2.00	0.60
1:F:169:ALA:HA	1:F:172:ARG:CZ	2.30	0.60
1:C:88:VAL:HB	1:C:224:MET:HE3	1.83	0.60
1:A:199:ILE:C	1:A:200:LEU:HG	2.27	0.60
1:B:36:ASN:N	1:B:37:PRO:HD3	2.15	0.60
1:B:74:VAL:O	1:B:78:VAL:HG23	2.01	0.60
3:B:1237:GMP:N3	3:B:1237:GMP:H2'	2.15	0.60
1:D:16:ARG:HB3	1:D:232:VAL:HB	1.84	0.60
1:E:72:ILE:O	1:E:76:GLU:HG3	2.02	0.60
1:B:35:GLN:HA	4:B:2007:HOH:O	2.02	0.60
1:C:213:PRO:HB2	1:C:216:VAL:HG23	1.83	0.59
1:F:212:ALA:HB1	1:F:213:PRO:HD2	1.84	0.59
3:E:1237:GMP:N3	3:E:1237:GMP:H2'	2.17	0.59
1:D:199:ILE:HG23	1:D:224:MET:HG3	1.83	0.59
1:B:19:LEU:HD21	1:B:77:LEU:CD1	2.33	0.59
1:F:56:VAL:HG11	1:F:233:LEU:HD22	1.83	0.59
1:A:219:GLU:OE2	1:A:219:GLU:HA	2.03	0.59
1:D:130:ASP:OD1	1:D:195:ARG:HG2	2.03	0.59
1:E:230:GLU:O	1:E:233:LEU:HB2	2.03	0.58
1:E:168:ARG:O	1:E:172:ARG:HG3	2.04	0.58
1:D:219:GLU:O	1:D:223:ARG:HG3	2.03	0.58
4:B:2033:HOH:O	1:C:139:ARG:HD3	2.02	0.58
1:D:36:ASN:HD22	1:D:36:ASN:H	1.51	0.58
1:B:138:TRP:CE2	1:B:150:VAL:HG21	2.39	0.57
1:D:36:ASN:N	1:D:36:ASN:ND2	2.49	0.57
1:F:88:VAL:HB	1:F:224:MET:HE3	1.85	0.57
1:B:203:SER:O	1:B:204:ASN:HB3	2.05	0.57
1:C:151:GLY:HA3	1:C:176:LEU:HD11	1.86	0.57
1:C:27:GLU:HG2	1:C:31:LYS:HE2	1.87	0.57
1:C:25:ARG:HD3	1:C:221:VAL:HG13	1.87	0.57
1:B:42:ASP:HA	1:B:47:TRP:HD1	1.69	0.57
1:D:82:ALA:HB1	1:D:85:LEU:HD11	1.87	0.57
1:B:83:ARG:HB2	4:B:2018:HOH:O	2.04	0.57
1:E:107:GLN:O	1:E:129:PRO:HD2	2.05	0.57
1:A:98:LEU:HD13	1:A:149:ARG:NH1	2.19	0.57
1:F:36:ASN:N	1:F:37:PRO:HD3	2.20	0.57
1:B:19:LEU:HD11	1:B:85:LEU:HB3	1.87	0.56
1:A:102:GLU:O	1:A:202:VAL:HG23	2.05	0.56
1:D:67:THR:N	1:D:68:PRO:CD	2.68	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:30:ALA:O	1:E:34:LEU:HB2	2.06	0.56
1:E:99:ALA:HA	1:E:205:ARG:NH2	2.20	0.56
1:E:230:GLU:OE2	1:E:230:GLU:HA	2.04	0.56
1:D:94:ALA:HB2	1:D:178:PHE:CE1	2.41	0.56
1:F:214:PRO:HG2	1:F:215:GLU:OE1	2.05	0.56
1:B:123:ARG:HH21	1:C:173:TYR:HE2	1.53	0.56
1:C:61:GLN:HE21	1:C:62:THR:H	1.52	0.56
1:C:115:THR:OG1	1:D:157:ASP:HB3	2.05	0.56
1:C:98:LEU:HD21	1:C:149:ARG:HG3	1.88	0.56
1:D:63:THR:HA	1:D:73:VAL:HG21	1.88	0.56
1:C:140:ARG:HA	1:C:140:ARG:NE	2.20	0.56
1:C:230:GLU:O	1:C:234:GLU:HG3	2.06	0.55
1:A:113:ASP:OD1	1:A:116:THR:HG23	2.06	0.55
1:D:180:MET:SD	3:D:1237:GMP:H3'	2.46	0.55
1:E:146:TYR:CG	1:E:223:ARG:HD3	2.41	0.55
1:C:203:SER:HB2	1:C:217:LEU:HD22	1.88	0.55
1:D:67:THR:HG22	1:D:181:GLU:HA	1.87	0.55
1:C:16:ARG:HH21	1:C:16:ARG:HG3	1.71	0.55
1:D:98:LEU:HD11	1:D:176:LEU:HB3	1.88	0.55
1:E:210:GLU:O	1:E:211:LEU:HD23	2.06	0.55
1:C:39:ARG:HD3	1:C:49:TYR:CE2	2.41	0.55
1:C:61:GLN:HE21	1:C:62:THR:N	2.05	0.55
1:D:3:PRO:CG	1:D:80:LEU:HD21	2.37	0.54
1:A:163:THR:HG22	4:A:2049:HOH:O	2.07	0.54
1:E:126:ALA:O	1:E:128:VAL:HG13	2.07	0.54
1:C:85:LEU:HD12	1:C:85:LEU:H	1.71	0.54
1:D:46:LEU:HA	1:D:62:THR:HB	1.90	0.54
1:F:102:GLU:O	1:F:202:VAL:HG23	2.07	0.54
1:B:229:LEU:O	1:B:233:LEU:HD13	2.07	0.54
1:C:3:PRO:HB3	1:C:76:GLU:HG2	1.89	0.54
1:A:63:THR:HA	1:A:73:VAL:HG21	1.89	0.54
1:D:98:LEU:HD13	1:D:104:ILE:HD11	1.90	0.54
3:D:1237:GMP:H2'	3:D:1237:GMP:N3	2.22	0.54
1:E:133:VAL:CG2	1:E:195:ARG:HB3	2.37	0.54
1:B:139:ARG:NH2	4:B:2033:HOH:O	2.33	0.54
1:D:58:VAL:HG11	1:D:229:LEU:CD2	2.38	0.54
1:C:158:ALA:HB1	1:C:161:ALA:HB2	1.90	0.53
1:D:3:PRO:HG2	1:D:80:LEU:HD21	1.88	0.53
1:D:217:LEU:C	1:D:217:LEU:HD23	2.33	0.53
1:F:234:GLU:HA	1:F:234:GLU:OE1	2.08	0.53
1:A:98:LEU:CD1	1:A:149:ARG:HG3	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:179:GLU:CD	1:D:181:GLU:H	2.15	0.53
1:A:110:VAL:HG13	1:F:125:TYR:HA	1.90	0.53
1:A:213:PRO:HG2	1:A:216:VAL:HG21	1.90	0.53
1:B:43:HIS:CE1	1:B:44:ARG:HG3	2.43	0.53
3:A:1237:GMP:N3	3:A:1237:GMP:H2'	2.22	0.53
1:C:179:GLU:CD	1:C:182:ALA:H	2.17	0.53
1:C:213:PRO:O	1:C:215:GLU:N	2.41	0.53
1:B:200:LEU:N	1:B:200:LEU:HD22	2.23	0.53
1:D:6:VAL:O	1:D:8:ALA:N	2.41	0.53
1:D:203:SER:HB2	1:D:211:LEU:HD22	1.91	0.53
1:C:17:VAL:HG22	1:C:59:SER:HB2	1.91	0.53
1:B:192:ARG:O	1:B:194:VAL:HG13	2.09	0.53
1:C:213:PRO:HB2	1:C:216:VAL:CG2	2.39	0.53
1:F:215:GLU:OE1	1:F:215:GLU:N	2.39	0.53
1:B:155:SER:HA	1:B:179:GLU:O	2.09	0.52
1:A:3:PRO:HB3	1:A:76:GLU:HG2	1.90	0.52
1:C:146:TYR:CD1	1:C:223:ARG:HB3	2.43	0.52
1:E:22:ASP:HB2	1:F:44:ARG:NH1	2.25	0.52
1:C:34:LEU:O	1:C:37:PRO:HD3	2.10	0.52
1:D:125:TYR:HA	1:E:110:VAL:HG13	1.92	0.52
1:F:85:LEU:O	1:F:196:THR:HA	2.09	0.52
1:E:133:VAL:HG21	1:E:195:ARG:HB3	1.92	0.52
1:F:36:ASN:ND2	4:F:2011:HOH:O	2.38	0.52
1:C:213:PRO:O	1:C:216:VAL:HG23	2.10	0.52
1:E:88:VAL:HG23	1:E:88:VAL:O	2.11	0.51
1:A:34:LEU:HD11	1:A:60:VAL:HG23	1.92	0.51
1:C:66:GLY:C	1:C:68:PRO:HD2	2.35	0.51
1:C:70:ALA:O	1:C:74:VAL:HG23	2.09	0.51
3:F:1237:GMP:H2'	3:F:1237:GMP:N3	2.25	0.51
1:E:221:VAL:O	1:E:224:MET:HB3	2.10	0.51
1:F:98:LEU:HD11	1:F:176:LEU:HB3	1.93	0.51
1:A:94:ALA:O	1:A:168:ARG:HG2	2.11	0.51
1:A:199:ILE:O	1:A:200:LEU:HG	2.11	0.51
1:D:140:ARG:O	1:D:142:GLU:N	2.43	0.51
1:B:146:TYR:CE1	1:B:223:ARG:HG2	2.46	0.51
1:B:36:ASN:N	1:B:36:ASN:HD22	2.08	0.51
1:F:199:ILE:HG12	1:F:224:MET:HG3	1.92	0.51
1:F:88:VAL:HG23	1:F:88:VAL:O	2.12	0.50
1:B:19:LEU:HD21	1:B:77:LEU:HD11	1.92	0.50
1:D:95:SER:HB2	1:D:97:ASP:OD2	2.12	0.50
1:F:52:LEU:HD22	1:F:55:GLY:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:THR:HG23	1:A:59:SER:OG	2.11	0.50
1:C:88:VAL:HG23	1:C:88:VAL:O	2.11	0.50
1:C:180:MET:SD	3:C:1237:GMP:H3'	2.52	0.50
1:E:9:HIS:O	1:E:12:ASP:HB2	2.12	0.50
1:B:53:TYR:HB3	1:B:58:VAL:HG21	1.93	0.50
1:B:181:GLU:O	1:B:182:ALA:C	2.55	0.50
1:D:98:LEU:HD23	1:D:149:ARG:NH1	2.26	0.50
1:D:140:ARG:C	1:D:142:GLU:N	2.70	0.50
1:A:65:MET:O	4:A:2021:HOH:O	2.20	0.50
1:A:135:ARG:O	1:A:139:ARG:HG3	2.12	0.50
1:A:180:MET:HE3	1:B:68:PRO:HB3	1.94	0.50
1:C:191:MET:HE2	1:C:192:ARG:NH2	2.20	0.50
1:E:205:ARG:O	1:E:206:ILE:C	2.55	0.50
1:B:85:LEU:HD12	1:B:85:LEU:N	2.27	0.49
1:E:135:ARG:HD3	4:E:2025:HOH:O	2.11	0.49
1:C:213:PRO:O	1:C:214:PRO:C	2.55	0.49
1:D:27:GLU:HG2	1:D:31:LYS:HE3	1.95	0.49
1:D:139:ARG:HE	1:E:139:ARG:NH2	2.09	0.49
1:C:146:TYR:HD1	1:C:223:ARG:HB3	1.76	0.49
1:D:130:ASP:OD2	1:D:195:ARG:NH1	2.44	0.49
1:A:168:ARG:HG3	1:A:168:ARG:HH11	1.76	0.49
1:E:63:THR:HA	1:E:73:VAL:HG21	1.95	0.49
1:A:67:THR:N	1:A:68:PRO:CD	2.76	0.49
1:B:36:ASN:N	1:B:36:ASN:ND2	2.59	0.49
1:B:140:ARG:O	1:B:144:LEU:HG	2.13	0.49
1:C:16:ARG:HG3	1:C:16:ARG:NH2	2.28	0.49
1:E:51:GLY:O	1:E:57:PRO:HA	2.12	0.49
1:F:98:LEU:CD2	1:F:149:ARG:HG3	2.39	0.49
1:F:203:SER:O	1:F:204:ASN:HB3	2.12	0.49
1:D:162:THR:HA	1:D:166:GLU:OE1	2.13	0.49
1:C:66:GLY:C	1:C:68:PRO:CD	2.86	0.49
1:B:219:GLU:OE2	1:B:219:GLU:HA	2.13	0.49
1:E:18:LEU:HD11	1:E:229:LEU:HG	1.94	0.49
1:E:79:ARG:O	1:E:79:ARG:NH1	2.45	0.49
1:C:68:PRO:HG2	1:D:68:PRO:HG2	1.94	0.48
1:B:135:ARG:NH1	1:C:138:TRP:CZ2	2.82	0.48
1:C:16:ARG:NE	1:C:233:LEU:HA	2.28	0.48
1:D:98:LEU:HD21	1:D:149:ARG:HG3	1.94	0.48
1:C:199:ILE:HG12	1:C:224:MET:HG3	1.94	0.48
1:D:140:ARG:C	1:D:142:GLU:H	2.22	0.48
1:B:18:LEU:HD11	1:B:229:LEU:HG	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:ALA:HB1	4:C:2001:HOH:O	2.13	0.48
1:A:6:VAL:O	1:A:8:ALA:N	2.46	0.48
1:B:8:ALA:HB3	1:B:80:LEU:HD13	1.96	0.48
1:C:129:PRO:CG	1:C:134:PHE:HB2	2.44	0.48
1:E:35:GLN:O	1:E:36:ASN:C	2.57	0.48
1:A:6:VAL:O	1:A:6:VAL:HG12	2.13	0.48
1:C:6:VAL:HG12	1:C:6:VAL:O	2.14	0.48
1:D:139:ARG:NE	1:E:139:ARG:HH22	2.10	0.48
1:A:44:ARG:HA	1:B:23:PRO:HD2	1.95	0.47
1:C:229:LEU:O	1:C:233:LEU:HD13	2.14	0.47
1:D:16:ARG:HD3	1:D:232:VAL:O	2.14	0.47
1:F:163:THR:OG1	1:F:166:GLU:HG3	2.14	0.47
1:E:9:HIS:CD2	1:E:11:GLY:H	2.32	0.47
1:D:72:ILE:O	1:D:76:GLU:HG3	2.14	0.47
1:D:178:PHE:CD1	1:D:206:ILE:CD1	2.95	0.47
1:D:90:THR:HA	1:D:201:ALA:O	2.13	0.47
1:A:120:LEU:O	1:A:121:GLU:C	2.57	0.47
1:A:166:GLU:O	1:A:169:ALA:HB3	2.15	0.47
1:C:16:ARG:HD3	1:C:232:VAL:C	2.39	0.47
1:B:228:ALA:O	1:B:232:VAL:HG22	2.15	0.47
1:C:61:GLN:NE2	1:C:62:THR:N	2.63	0.47
1:C:88:VAL:CB	1:C:224:MET:HE3	2.45	0.47
1:D:94:ALA:HB2	1:D:178:PHE:HE1	1.77	0.47
1:D:116:THR:HG22	1:D:184:ALA:HA	1.97	0.47
1:F:85:LEU:HD12	1:F:85:LEU:N	2.30	0.47
1:A:131:PRO:HG2	4:A:2042:HOH:O	2.14	0.47
1:D:205:ARG:O	1:D:206:ILE:C	2.58	0.47
1:E:208:ASP:HA	1:E:209:PRO:HD3	1.73	0.47
1:F:9:HIS:HB3	1:F:10:PRO:HD2	1.96	0.47
1:D:39:ARG:HA	1:D:49:TYR:HA	1.97	0.47
1:E:146:TYR:CD1	1:E:223:ARG:HB3	2.49	0.47
1:A:103:LEU:HA	1:A:200:LEU:O	2.14	0.46
1:C:28:TRP:O	1:C:32:THR:OG1	2.18	0.46
1:B:23:PRO:HA	1:B:62:THR:OG1	2.15	0.46
1:D:38:ARG:HG2	1:D:38:ARG:HH11	1.80	0.46
1:A:161:ALA:HB3	1:B:118:GLN:HE21	1.79	0.46
1:B:135:ARG:HD2	1:C:138:TRP:CH2	2.51	0.46
1:D:229:LEU:O	1:D:233:LEU:HD13	2.16	0.46
1:B:9:HIS:HB3	1:B:10:PRO:CD	2.42	0.46
1:B:140:ARG:HB2	1:B:227:VAL:HG13	1.97	0.46
1:C:67:THR:N	1:C:68:PRO:CD	2.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:228:ALA:O	1:E:232:VAL:HG13	2.16	0.46
1:C:229:LEU:O	1:C:232:VAL:HG23	2.15	0.46
1:C:146:TYR:N	1:C:146:TYR:CD2	2.82	0.46
1:A:61:GLN:HG2	1:A:77:LEU:HD11	1.99	0.45
1:B:135:ARG:NH1	1:C:138:TRP:CE2	2.84	0.45
1:E:87:ARG:NH2	1:E:89:GLY:HA3	2.31	0.45
1:F:205:ARG:NH2	4:F:2056:HOH:O	2.47	0.45
1:F:213:PRO:HA	1:F:214:PRO:HD2	1.81	0.45
1:A:210:GLU:C	1:A:211:LEU:HD23	2.41	0.45
1:D:221:VAL:HG13	1:D:224:MET:HE2	1.99	0.45
1:E:25:ARG:O	1:E:29:ILE:HG13	2.16	0.45
1:A:217:LEU:O	1:A:221:VAL:HG23	2.16	0.45
1:B:77:LEU:HB3	1:B:82:ALA:HB2	1.98	0.45
1:C:98:LEU:HD11	1:C:176:LEU:HB3	1.98	0.45
1:C:155:SER:OG	1:C:179:GLU:OE1	2.26	0.45
1:E:117:ARG:NH1	4:E:2018:HOH:O	2.49	0.45
1:E:180:MET:O	1:E:181:GLU:HG3	2.17	0.45
1:C:199:ILE:HG23	1:C:224:MET:CG	2.47	0.45
1:D:74:VAL:O	1:D:78:VAL:HG23	2.17	0.45
1:E:113:ASP:OD2	1:E:116:THR:HG23	2.16	0.45
1:F:17:VAL:HA	1:F:59:SER:O	2.17	0.45
1:B:16:ARG:HH21	1:B:16:ARG:HG3	1.82	0.45
1:B:139:ARG:HB3	1:B:140:ARG:HH21	1.81	0.45
1:D:21:GLY:HA3	2:D:1236:SO4:O2	2.16	0.45
1:A:112:LEU:HB2	1:A:155:SER:O	2.16	0.45
1:B:144:LEU:HD13	1:B:226:GLU:CD	2.42	0.45
1:B:159:PHE:C	1:B:159:PHE:CD2	2.94	0.45
1:D:90:THR:O	3:D:1237:GMP:H1'	2.16	0.45
1:E:19:LEU:HB2	1:E:87:ARG:HA	1.99	0.45
1:A:168:ARG:HG3	1:A:168:ARG:NH1	2.32	0.45
1:B:171:ALA:C	1:B:173:TYR:H	2.24	0.45
1:C:110:VAL:HG12	1:C:112:LEU:HG	1.98	0.45
1:D:18:LEU:HD11	1:D:229:LEU:HG	1.98	0.45
1:E:135:ARG:HG3	1:E:139:ARG:NH1	2.32	0.45
1:A:56:VAL:HG13	1:A:57:PRO:HD2	1.97	0.45
1:A:203:SER:O	1:A:204:ASN:HB3	2.17	0.45
1:C:235:VAL:HG23	1:C:235:VAL:OXT	2.16	0.45
1:D:188:LEU:HA	1:D:191:MET:HB2	1.99	0.44
1:A:51:GLY:O	1:A:57:PRO:HA	2.17	0.44
1:C:61:GLN:HB3	1:C:77:LEU:HD11	1.99	0.44
1:C:203:SER:HA	1:C:217:LEU:HD13	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:229:LEU:O	1:C:232:VAL:CG2	2.65	0.44
1:D:36:ASN:N	1:D:37:PRO:CD	2.80	0.44
1:D:209:PRO:C	1:D:210:GLU:HG3	2.43	0.44
1:E:180:MET:SD	3:E:1237:GMP:H3'	2.57	0.44
1:D:92:GLY:HA3	3:D:1237:GMP:O6	2.18	0.44
1:D:139:ARG:NE	1:E:139:ARG:NH2	2.66	0.44
1:B:209:PRO:HG2	1:B:210:GLU:H	1.80	0.44
1:B:213:PRO:HA	1:B:214:PRO:HD2	1.78	0.44
1:E:9:HIS:HD2	1:E:11:GLY:H	1.65	0.44
1:E:230:GLU:O	1:E:231:ALA:C	2.60	0.44
1:C:5:HIS:CD2	1:C:72:ILE:HG21	2.53	0.44
1:A:209:PRO:HD2	4:A:2054:HOH:O	2.17	0.44
1:C:16:ARG:HE	1:C:233:LEU:HA	1.81	0.44
1:C:35:GLN:O	1:C:36:ASN:C	2.61	0.44
1:C:52:LEU:HD13	1:C:55:GLY:O	2.18	0.44
1:E:215:GLU:CB	4:E:2036:HOH:O	2.66	0.44
1:A:164:PRO:HG3	1:A:207:GLY:HA2	1.98	0.44
1:B:90:THR:O	3:B:1237:GMP:H1'	2.18	0.44
1:C:231:ALA:O	1:C:234:GLU:HB2	2.17	0.44
1:D:85:LEU:N	1:D:85:LEU:CD1	2.80	0.44
1:B:41:ASN:O	1:B:47:TRP:HA	2.17	0.44
1:B:63:THR:HA	1:B:73:VAL:HG21	1.99	0.44
1:C:95:SER:HB2	1:C:97:ASP:OD2	2.17	0.44
1:D:94:ALA:CB	1:D:178:PHE:HE1	2.31	0.44
1:F:136:ALA:O	1:F:140:ARG:HG2	2.17	0.44
1:F:149:ARG:NH1	1:F:149:ARG:HB3	2.33	0.44
1:C:137:LEU:HD23	1:C:231:ALA:CB	2.47	0.43
1:B:229:LEU:O	1:B:232:VAL:HG23	2.18	0.43
1:F:149:ARG:HB3	1:F:149:ARG:HH11	1.83	0.43
1:B:36:ASN:HD22	1:B:36:ASN:H	1.66	0.43
1:B:171:ALA:HB3	4:B:2038:HOH:O	2.18	0.43
1:B:226:GLU:O	1:B:227:VAL:C	2.60	0.43
1:C:58:VAL:HG12	1:C:59:SER:N	2.32	0.43
1:C:135:ARG:HD3	1:C:139:ARG:NH2	2.34	0.43
1:D:138:TRP:CE2	1:E:135:ARG:NH2	2.86	0.43
1:E:113:ASP:OD1	1:E:116:THR:HG23	2.18	0.43
1:E:212:ALA:CB	1:E:213:PRO:CD	2.89	0.43
1:C:190:ARG:HG2	4:C:2017:HOH:O	2.18	0.43
1:E:25:ARG:HH12	1:E:89:GLY:HA2	1.83	0.43
1:B:42:ASP:HA	1:B:47:TRP:CD1	2.52	0.43
1:D:110:VAL:HG13	1:E:125:TYR:HA	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:204:ASN:ND2	1:E:205:ARG:N	2.63	0.43
1:A:219:GLU:O	1:A:223:ARG:HG3	2.17	0.43
1:C:52:LEU:HD22	1:C:55:GLY:O	2.18	0.43
1:C:65:MET:HA	1:C:181:GLU:OE1	2.18	0.43
1:C:20:PRO:HB3	1:C:88:VAL:HG23	2.00	0.43
1:E:37:PRO:C	1:E:38:ARG:HD3	2.44	0.43
1:B:168:ARG:HG3	4:B:2020:HOH:O	2.18	0.43
1:C:208:ASP:HA	1:C:209:PRO:HD3	1.70	0.43
1:D:107:GLN:HA	1:D:134:PHE:CE1	2.54	0.43
1:A:30:ALA:HA	1:A:34:LEU:HG	2.00	0.43
1:B:123:ARG:HD2	1:C:170:TRP:CZ2	2.53	0.43
1:C:130:ASP:OD2	1:C:132:GLU:HB2	2.19	0.43
1:F:17:VAL:HG22	1:F:59:SER:HB2	2.01	0.43
1:C:108:GLY:C	1:C:152:LEU:HD23	2.43	0.42
1:D:130:ASP:CG	1:D:195:ARG:HH11	2.27	0.42
1:C:112:LEU:HD21	1:C:170:TRP:CZ2	2.54	0.42
1:D:106:ALA:HB3	1:D:198:ALA:HB3	2.01	0.42
1:D:140:ARG:HA	1:D:140:ARG:NE	2.34	0.42
1:B:53:TYR:CE2	1:B:229:LEU:HB3	2.54	0.42
1:B:119:TYR:CZ	1:B:188:LEU:HD21	2.54	0.42
1:B:130:ASP:OD2	1:B:133:VAL:HG23	2.18	0.42
1:D:38:ARG:HG2	1:D:38:ARG:NH1	2.34	0.42
1:D:67:THR:N	1:D:68:PRO:HD3	2.35	0.42
1:F:91:ALA:HB2	1:F:179:GLU:HB3	2.00	0.42
1:F:120:LEU:O	1:F:121:GLU:HB2	2.20	0.42
1:C:188:LEU:HA	1:C:191:MET:HB2	2.01	0.42
1:A:85:LEU:O	1:A:196:THR:HA	2.19	0.42
1:C:17:VAL:HG13	1:C:59:SER:HB3	2.02	0.42
1:D:146:TYR:CE1	1:D:223:ARG:HG2	2.55	0.42
1:E:56:VAL:HA	1:E:57:PRO:HD3	1.82	0.42
1:E:199:ILE:HG23	1:E:199:ILE:O	2.19	0.42
1:A:223:ARG:O	1:A:227:VAL:HG23	2.19	0.42
1:C:85:LEU:O	1:C:196:THR:HA	2.20	0.42
1:C:199:ILE:C	1:C:200:LEU:HG	2.44	0.42
1:C:129:PRO:HG2	1:C:134:PHE:HB2	2.01	0.42
1:D:107:GLN:OE1	1:E:131:PRO:HG3	2.19	0.42
1:A:85:LEU:N	1:A:85:LEU:CD1	2.80	0.41
1:B:158:ALA:CB	1:B:161:ALA:HB2	2.50	0.41
1:E:98:LEU:HA	1:E:102:GLU:OE2	2.20	0.41
1:E:180:MET:HG3	3:E:1237:GMP:HN22	1.86	0.41
1:F:56:VAL:HG11	1:F:233:LEU:CD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:13:VAL:HG21	1:D:80:LEU:HD12	2.02	0.41
1:F:209:PRO:O	1:F:210:GLU:HG3	2.19	0.41
1:A:3:PRO:HB3	1:A:76:GLU:CG	2.51	0.41
1:D:213:PRO:HA	1:D:214:PRO:HD3	1.90	0.41
1:A:156:GLU:O	1:A:180:MET:HG2	2.21	0.41
1:C:158:ALA:CB	1:C:161:ALA:HB2	2.50	0.41
1:A:13:VAL:HG22	1:A:40:TYR:CZ	2.56	0.41
1:B:67:THR:N	1:B:68:PRO:CD	2.84	0.41
1:C:130:ASP:OD1	1:C:195:ARG:HG2	2.20	0.41
1:D:123:ARG:HD2	1:E:170:TRP:CZ2	2.56	0.41
1:E:146:TYR:CD2	1:E:223:ARG:HD3	2.55	0.41
1:E:213:PRO:HB2	1:E:216:VAL:HG23	2.02	0.41
1:F:2:SER:N	4:F:2001:HOH:O	2.54	0.41
1:F:67:THR:N	1:F:68:PRO:CD	2.84	0.41
1:C:97:ASP:OD2	1:C:97:ASP:N	2.50	0.41
1:D:6:VAL:O	1:D:6:VAL:HG12	2.21	0.41
1:D:123:ARG:HA	1:D:124:PRO:HD3	1.87	0.41
1:F:10:PRO:HA	4:F:2022:HOH:O	2.20	0.41
1:B:130:ASP:O	1:B:131:PRO:C	2.63	0.40
1:D:123:ARG:NH1	1:D:123:ARG:HG3	2.36	0.40
1:D:138:TRP:CE2	1:D:150:VAL:HG21	2.56	0.40
1:F:125:TYR:HD2	1:F:127:PRO:HG3	1.85	0.40
1:C:179:GLU:C	1:C:180:MET:HG3	2.46	0.40
1:E:85:LEU:N	1:E:85:LEU:HD12	2.35	0.40
1:E:98:LEU:HD23	1:E:102:GLU:CD	2.47	0.40
1:F:129:PRO:HB3	1:F:186:PHE:CZ	2.56	0.40
1:C:127:PRO:HA	4:C:2010:HOH:O	2.20	0.40
1:D:144:LEU:O	1:D:146:TYR:N	2.54	0.40
1:E:124:PRO:O	1:E:125:TYR:HB2	2.20	0.40
1:A:107:GLN:HA	1:A:134:PHE:CE1	2.56	0.40
1:B:96:SER:HB2	4:B:2043:HOH:O	2.21	0.40
1:C:3:PRO:HB3	1:C:76:GLU:CG	2.51	0.40
1:D:7:ARG:NE	1:D:43:HIS:CE1	2.90	0.40
1:F:3:PRO:HG2	1:F:80:LEU:HD11	2.04	0.40
1:F:205:ARG:NH1	1:F:205:ARG:HG2	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	232/235 (99%)	213 (92%)	16 (7%)	3 (1%)	9	14
1	B	232/235 (99%)	207 (89%)	20 (9%)	5 (2%)	5	6
1	C	232/235 (99%)	198 (85%)	27 (12%)	7 (3%)	3	3
1	D	232/235 (99%)	194 (84%)	32 (14%)	6 (3%)	4	4
1	E	232/235 (99%)	197 (85%)	32 (14%)	3 (1%)	9	14
1	F	232/235 (99%)	218 (94%)	13 (6%)	1 (0%)	30	43
All	All	1392/1410 (99%)	1227 (88%)	140 (10%)	25 (2%)	6	9

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	ARG
1	C	214	PRO
1	D	7	ARG
1	D	206	ILE
1	D	211	LEU
1	C	204	ASN
1	D	141	ALA
1	D	145	GLY
1	A	214	PRO
1	C	213	PRO
1	B	182	ALA
1	B	204	ASN
1	C	7	ARG
1	F	204	ASN
1	B	213	PRO
1	C	30	ALA
1	E	206	ILE
1	E	231	ALA
1	A	210	GLU
1	C	20	PRO

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Mol	Chain	Res	Type
1	B	206	ILE
1	B	193	GLY
1	C	153	VAL
1	D	213	PRO
1	E	212	ALA

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	179/180 (99%)	173 (97%)	6 (3%)	32	54
1	B	179/180 (99%)	169 (94%)	10 (6%)	19	33
1	C	179/180 (99%)	173 (97%)	6 (3%)	32	54
1	D	179/180 (99%)	172 (96%)	7 (4%)	28	48
1	E	179/180 (99%)	175 (98%)	4 (2%)	45	67
1	F	179/180 (99%)	176 (98%)	3 (2%)	53	74
All	All	1074/1080 (99%)	1038 (97%)	36 (3%)	32	54

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	85	LEU
1	A	88	VAL
1	A	95	SER
1	A	179	GLU
1	A	200	LEU
1	A	214	PRO
1	B	10	PRO
1	B	85	LEU
1	B	88	VAL
1	B	128	VAL
1	B	130	ASP
1	B	179	GLU
1	B	200	LEU

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Mol	Chain	Res	Type
1	B	213	PRO
1	B	217	LEU
1	B	232	VAL
1	C	59	SER
1	C	85	LEU
1	C	162	THR
1	C	214	PRO
1	C	215	GLU
1	C	216	VAL
1	D	22	ASP
1	D	36	ASN
1	D	88	VAL
1	D	95	SER
1	D	194	VAL
1	D	204	ASN
1	D	216	VAL
1	E	13	VAL
1	E	130	ASP
1	E	179	GLU
1	E	233	LEU
1	F	76	GLU
1	F	147	PRO
1	F	179	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	A	218	GLN
1	B	36	ASN
1	B	204	ASN
1	C	9	HIS
1	C	61	GLN
1	D	9	HIS
1	D	36	ASN
1	D	204	ASN
1	E	9	HIS
1	E	35	GLN
1	E	204	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

12 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	$\# Z > 2$
2	SO4	D	1236	-	4,4,4	0.40	0	6,6,6	0.22	0
2	SO4	E	1236	-	4,4,4	0.38	0	6,6,6	0.06	0
2	SO4	C	1236	-	4,4,4	0.39	0	6,6,6	0.14	0
3	GMP	A	1237	-	22,22,22	0.67	0	33,33,33	1.09	5 (15%)
3	GMP	D	1237	-	22,22,22	0.56	0	33,33,33	1.23	5 (15%)
3	GMP	F	1237	-	22,22,22	0.66	0	33,33,33	1.09	4 (12%)
2	SO4	A	1236	-	4,4,4	0.43	0	6,6,6	0.17	0
3	GMP	B	1237	-	22,22,22	0.76	0	33,33,33	1.02	4 (12%)
3	GMP	E	1237	-	22,22,22	0.63	0	33,33,33	1.09	4 (12%)
3	GMP	C	1237	-	22,22,22	0.65	0	33,33,33	1.12	4 (12%)
2	SO4	B	1236	-	4,4,4	0.31	0	6,6,6	0.16	0
2	SO4	F	1236	-	4,4,4	0.30	0	6,6,6	0.11	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GMP	A	1237	-	-	2/6/22/22	0/3/3/3
3	GMP	D	1237	-	-	1/6/22/22	0/3/3/3
3	GMP	F	1237	-	-	2/6/22/22	0/3/3/3
3	GMP	B	1237	-	-	2/6/22/22	0/3/3/3
3	GMP	C	1237	-	-	2/6/22/22	0/3/3/3
3	GMP	E	1237	-	-	2/6/22/22	0/3/3/3

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1237	GMP	O4'-C1'-N9	3.25	115.73	108.36
3	D	1237	GMP	O4'-C1'-N9	2.99	115.13	108.36
3	D	1237	GMP	C2'-C3'-C4'	2.98	108.36	102.61
3	D	1237	GMP	O4'-C4'-C5'	2.91	115.36	109.22
3	F	1237	GMP	O4'-C4'-C5'	2.76	115.06	109.22
3	A	1237	GMP	O4'-C4'-C5'	2.69	114.91	109.22
3	C	1237	GMP	C2'-C3'-C4'	2.59	107.61	102.61
3	E	1237	GMP	O4'-C4'-C5'	2.53	114.57	109.22
3	D	1237	GMP	O3'-C3'-C4'	-2.50	103.90	111.08
3	E	1237	GMP	C2'-C3'-C4'	2.47	107.39	102.61
3	C	1237	GMP	O5'-C5'-C4'	-2.45	103.00	111.33
3	E	1237	GMP	O4'-C1'-N9	2.33	113.64	108.36
3	C	1237	GMP	O4'-C4'-C5'	2.30	114.09	109.22
3	B	1237	GMP	O5'-C5'-C4'	-2.28	103.56	111.33
3	B	1237	GMP	O4'-C4'-C5'	2.27	114.02	109.22
3	A	1237	GMP	O5'-C5'-C4'	-2.25	103.68	111.33
3	D	1237	GMP	O5'-C5'-C4'	-2.25	103.68	111.33
3	F	1237	GMP	O3'-C3'-C4'	-2.24	104.65	111.08
3	A	1237	GMP	C2'-C3'-C4'	2.21	106.87	102.61
3	B	1237	GMP	O4'-C1'-N9	2.16	113.26	108.36
3	F	1237	GMP	O5'-C5'-C4'	-2.14	104.06	111.33
3	B	1237	GMP	C2'-C3'-C4'	2.11	106.69	102.61
3	A	1237	GMP	O4'-C1'-N9	2.04	112.98	108.36
3	A	1237	GMP	O3'-C3'-C4'	-2.01	105.30	111.08
3	E	1237	GMP	O3'-C3'-C4'	-2.01	105.31	111.08
3	F	1237	GMP	C2'-C3'-C4'	2.01	106.49	102.61

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	1237	GMP	C2'-C1'-N9-C4
3	B	1237	GMP	C2'-C1'-N9-C4
3	A	1237	GMP	C2'-C1'-N9-C4
3	C	1237	GMP	C2'-C1'-N9-C4
3	F	1237	GMP	C2'-C1'-N9-C4
3	E	1237	GMP	C2'-C1'-N9-C8
3	D	1237	GMP	C2'-C1'-N9-C4
3	B	1237	GMP	C2'-C1'-N9-C8
3	A	1237	GMP	C2'-C1'-N9-C8
3	C	1237	GMP	C2'-C1'-N9-C8
3	F	1237	GMP	C2'-C1'-N9-C8

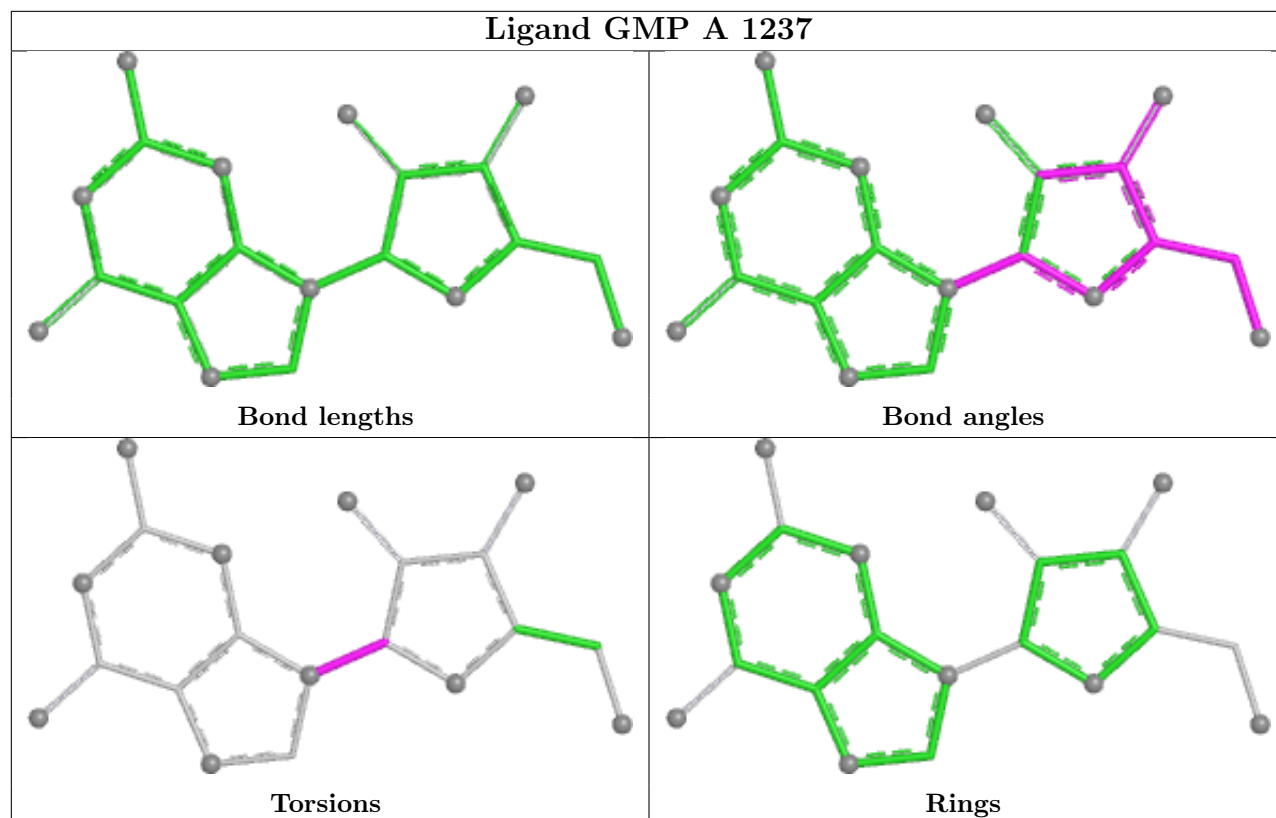
There are no ring outliers.

7 monomers are involved in 14 short contacts:

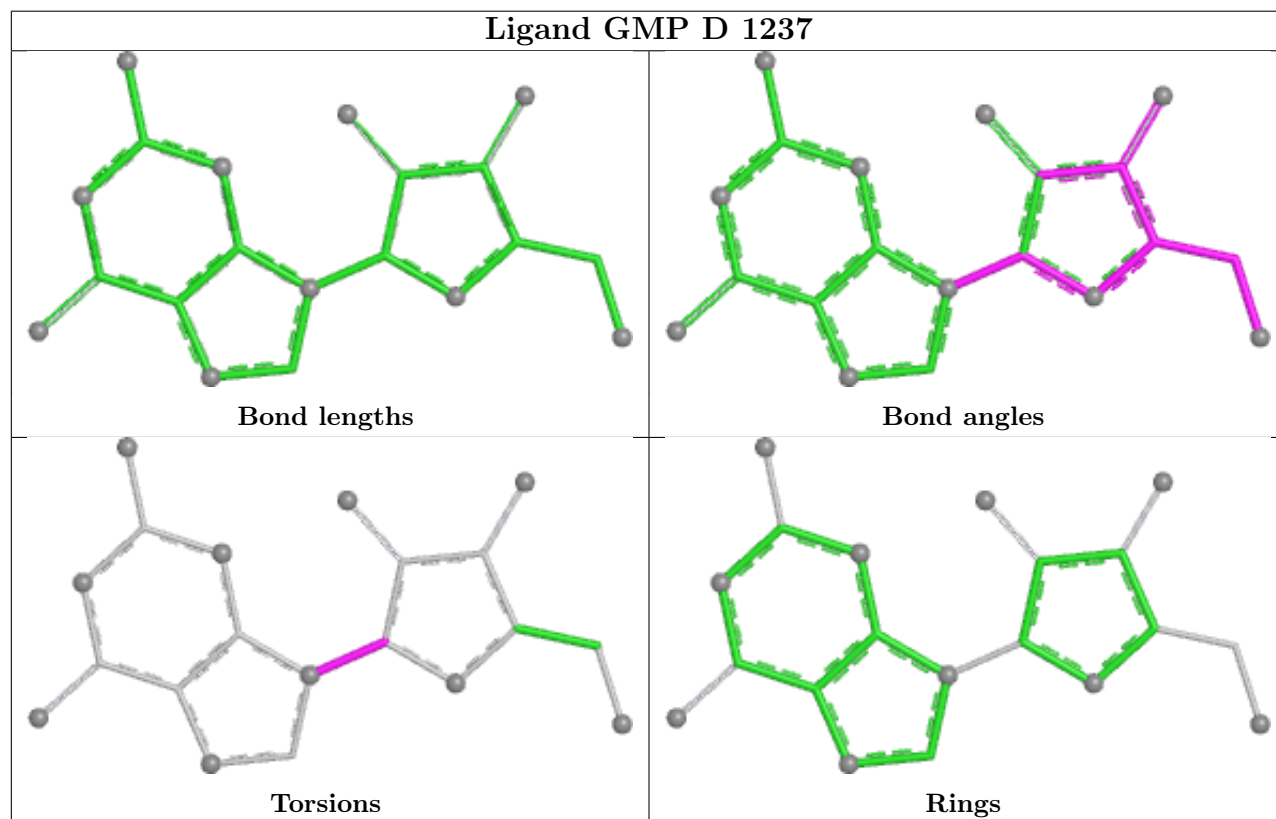
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1236	SO4	1	0
3	A	1237	GMP	1	0
3	D	1237	GMP	4	0
3	F	1237	GMP	1	0
3	B	1237	GMP	2	0
3	E	1237	GMP	3	0
3	C	1237	GMP	2	0

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.

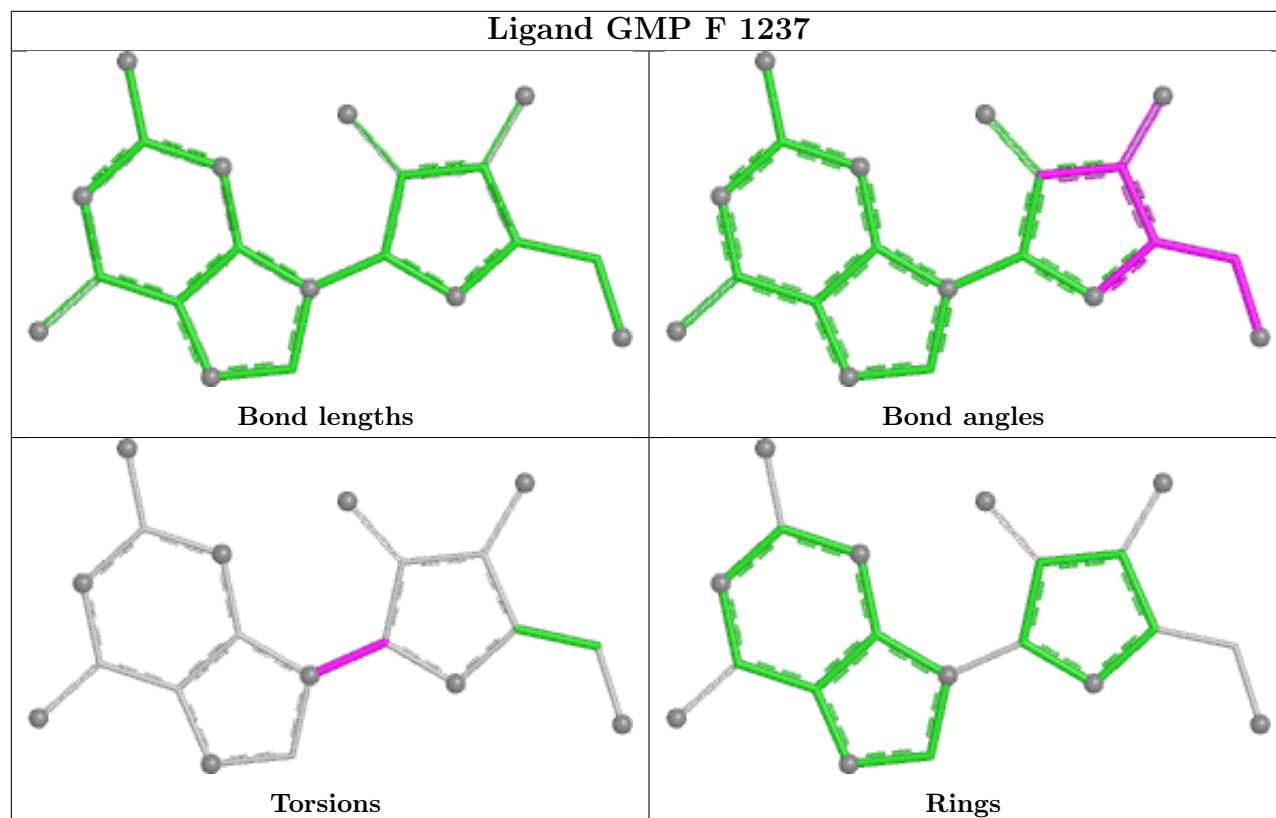
Ligand GMP A 1237



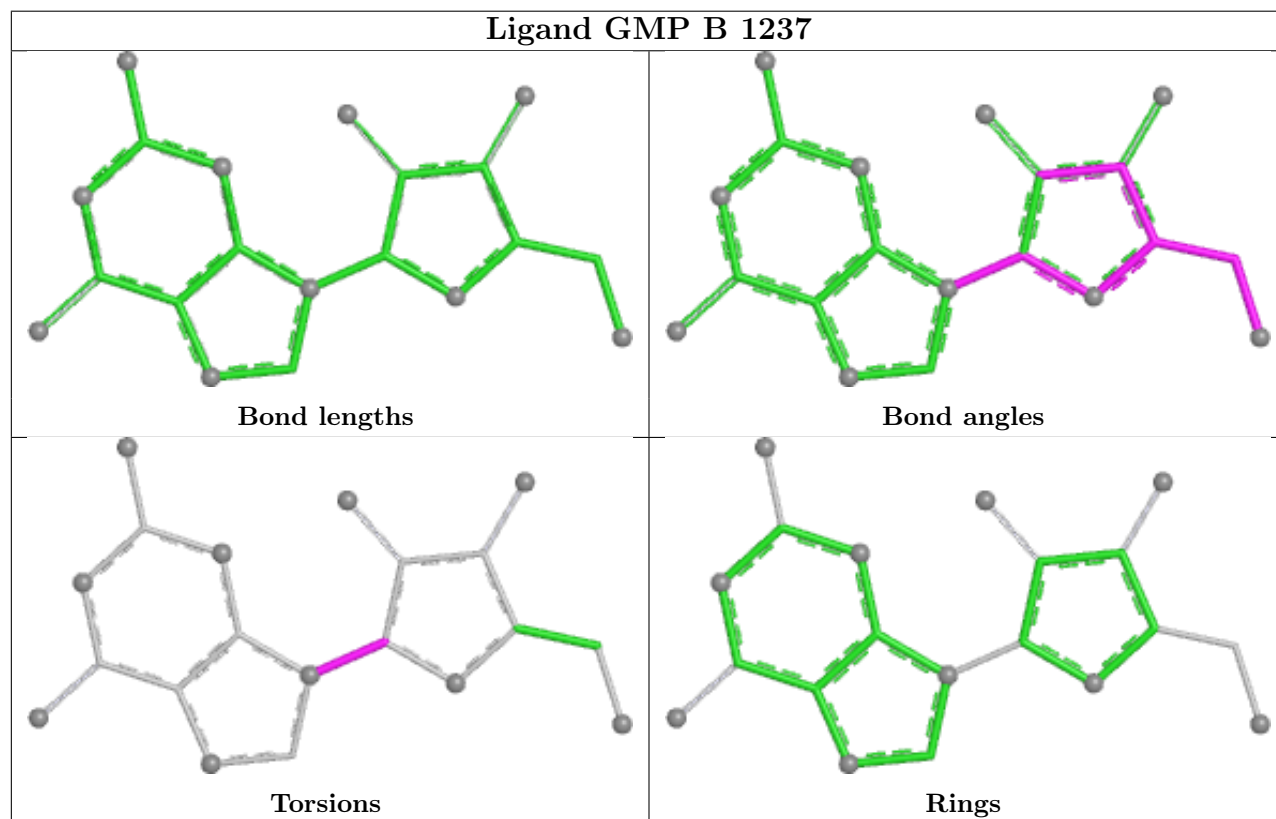
Ligand GMP D 1237



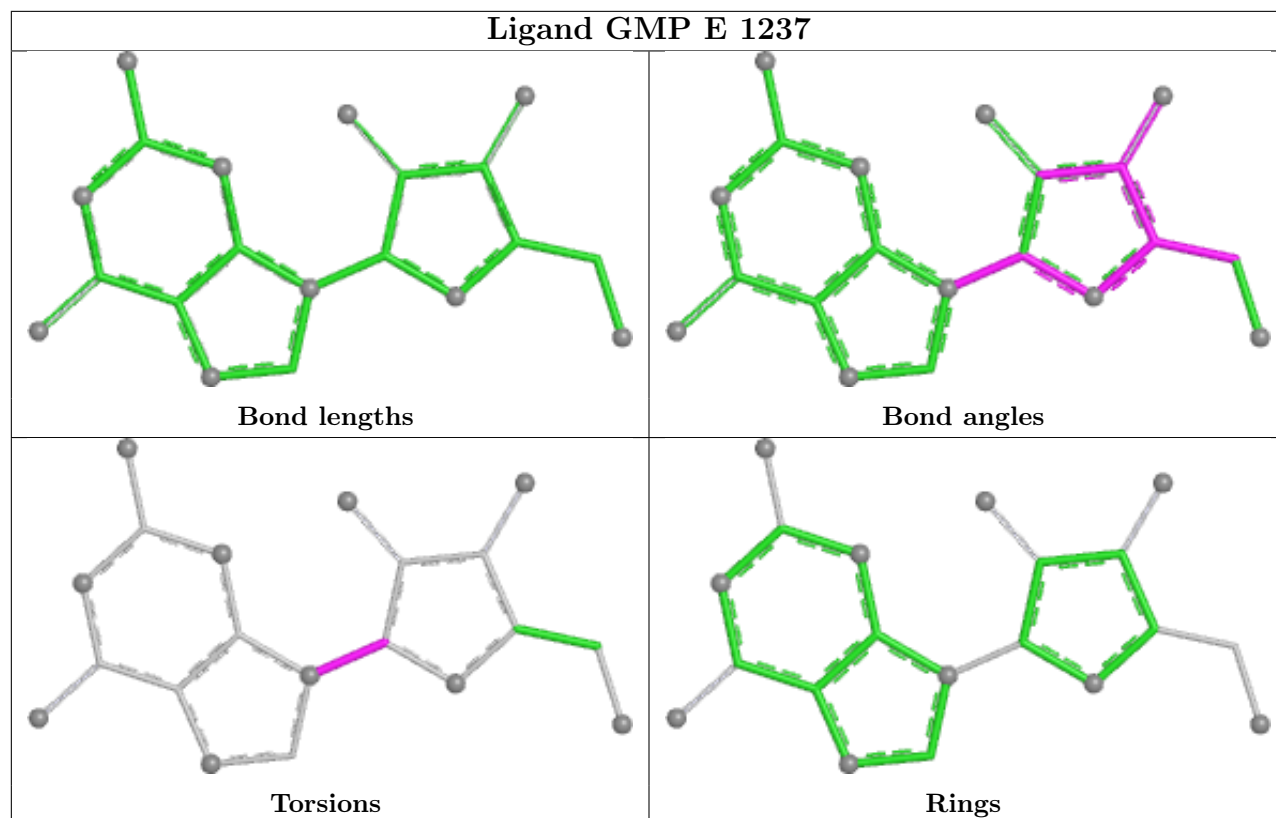
Ligand GMP F 1237



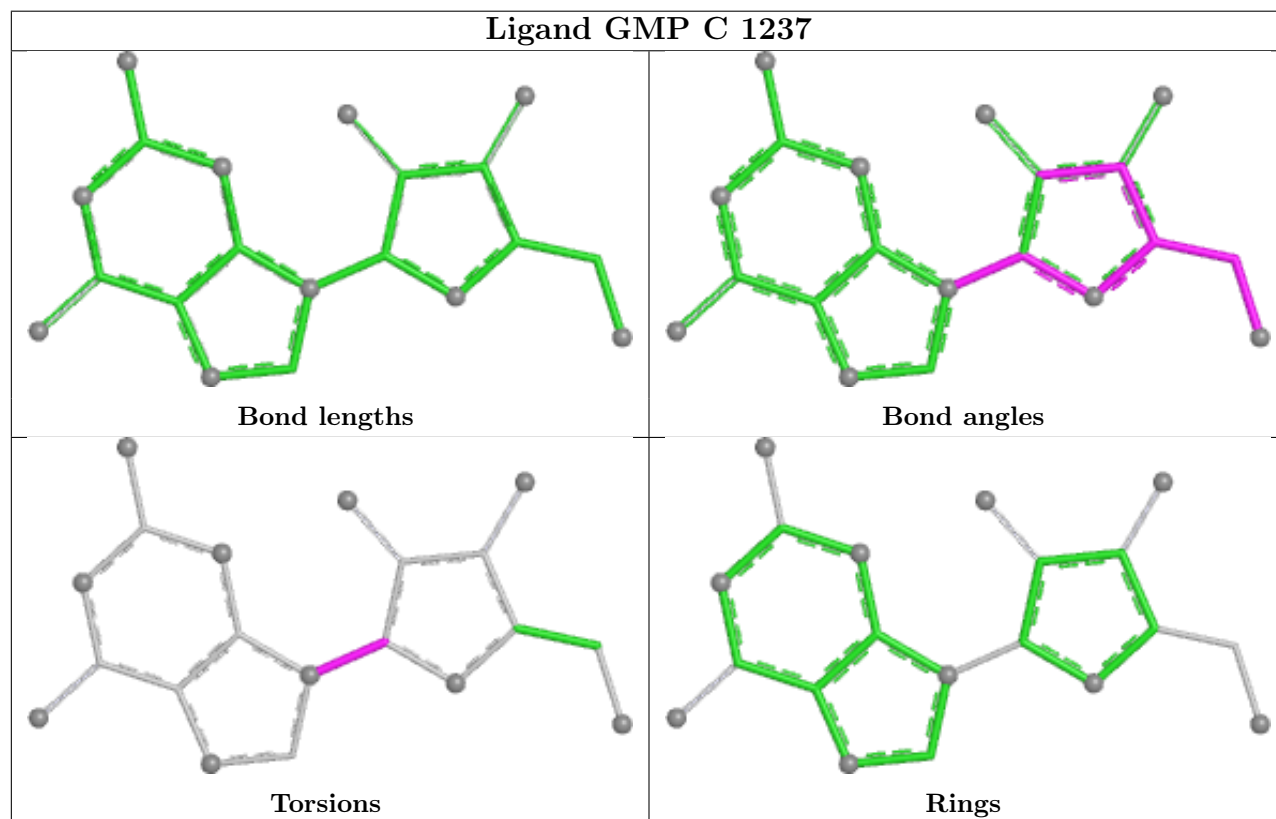
Ligand GMP B 1237



Ligand GMP E 1237



Ligand GMP C 1237



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	234/235 (99%)	-0.42	6 (2%) 57 53	3, 15, 45, 88	0
1	B	234/235 (99%)	-0.14	3 (1%) 75 71	10, 24, 53, 76	0
1	C	234/235 (99%)	0.17	8 (3%) 48 44	10, 31, 63, 83	0
1	D	234/235 (99%)	0.25	7 (2%) 52 48	11, 35, 61, 81	0
1	E	234/235 (99%)	-0.19	8 (3%) 48 44	3, 23, 56, 83	0
1	F	234/235 (99%)	-0.42	5 (2%) 63 59	3, 15, 44, 81	0
All	All	1404/1410 (99%)	-0.12	37 (2%) 57 53	3, 24, 57, 88	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	212	ALA	4.5
1	F	209	PRO	4.3
1	D	209	PRO	4.1
1	B	212	ALA	4.0
1	C	207	GLY	3.8
1	E	211	LEU	3.7
1	A	211	LEU	3.7
1	C	211	LEU	3.6
1	E	206	ILE	3.6
1	F	211	LEU	3.5
1	E	209	PRO	3.4
1	B	211	LEU	3.3
1	D	212	ALA	3.2
1	A	209	PRO	3.2
1	C	208	ASP	3.0
1	D	205	ARG	3.0
1	A	96	SER	2.6
1	D	213	PRO	2.6
1	C	212	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	210	GLU	2.5
1	E	210	GLU	2.5
1	E	207	GLY	2.5
1	F	208	ASP	2.5
1	C	98	LEU	2.4
1	C	217	LEU	2.4
1	D	211	LEU	2.4
1	E	213	PRO	2.3
1	D	206	ILE	2.3
1	C	209	PRO	2.3
1	F	206	ILE	2.2
1	C	216	VAL	2.2
1	A	207	GLY	2.2
1	D	38	ARG	2.2
1	B	209	PRO	2.1
1	E	212	ALA	2.1
1	E	217	LEU	2.1
1	A	215	GLU	2.0

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	GMP	C	1237	20/20	0.85	0.12	43,54,59,60	0
3	GMP	D	1237	20/20	0.87	0.12	32,37,42,46	0
3	GMP	F	1237	20/20	0.94	0.07	14,19,24,26	0
2	SO4	C	1236	5/5	0.95	0.08	30,35,40,40	0
3	GMP	A	1237	20/20	0.95	0.06	13,19,26,31	0

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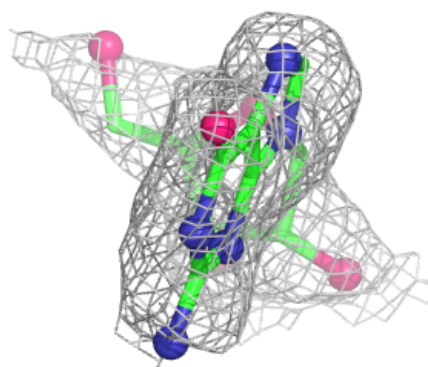
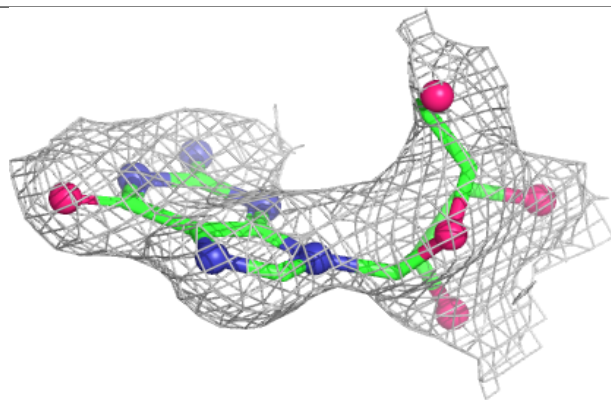
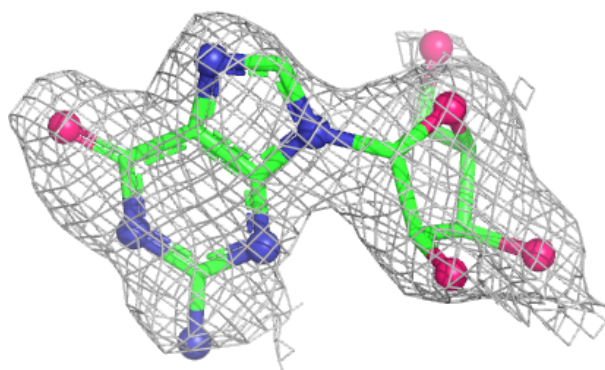
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GMP	B	1237	20/20	0.95	0.07	11,15,19,23	0
3	GMP	E	1237	20/20	0.96	0.06	12,17,20,21	0
2	SO4	D	1236	5/5	0.97	0.06	24,28,31,31	0
2	SO4	E	1236	5/5	0.99	0.05	25,26,30,31	0
2	SO4	F	1236	5/5	0.99	0.04	17,19,21,23	0
2	SO4	A	1236	5/5	0.99	0.05	13,14,16,18	0
2	SO4	B	1236	5/5	0.99	0.04	14,15,15,16	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.

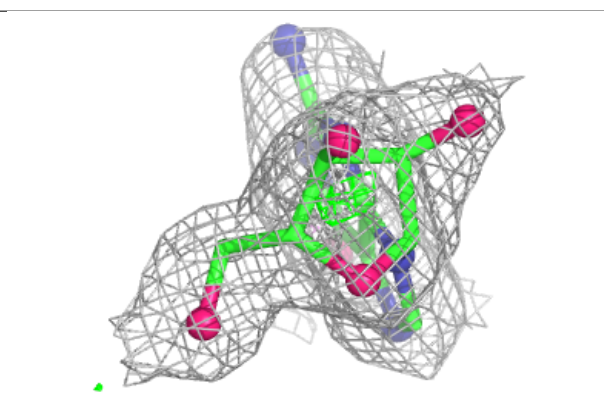
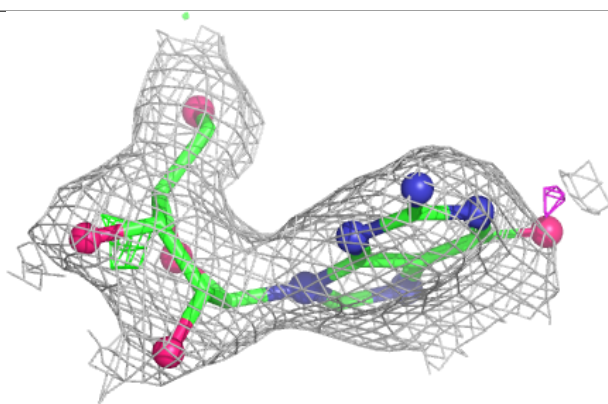
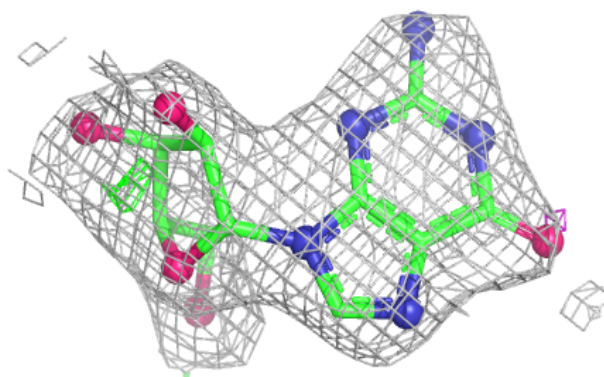
Electron density around GMP C 1237:

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

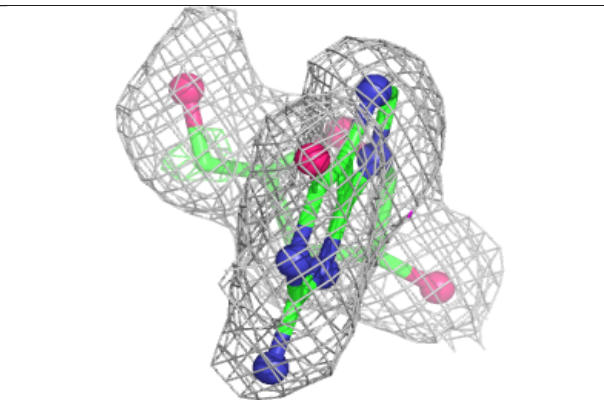
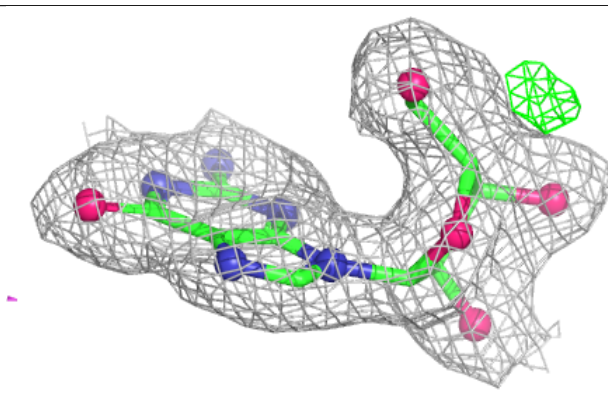
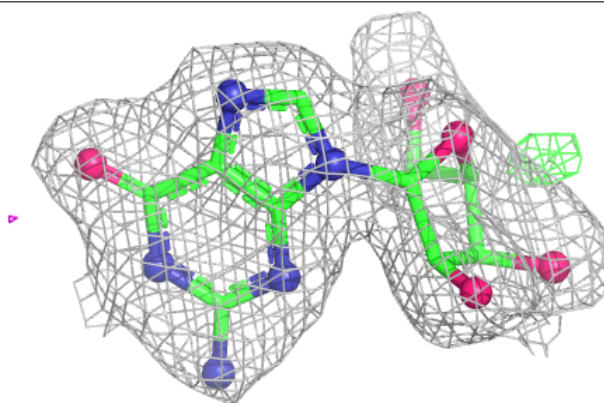


Electron density around GMP D 1237:

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

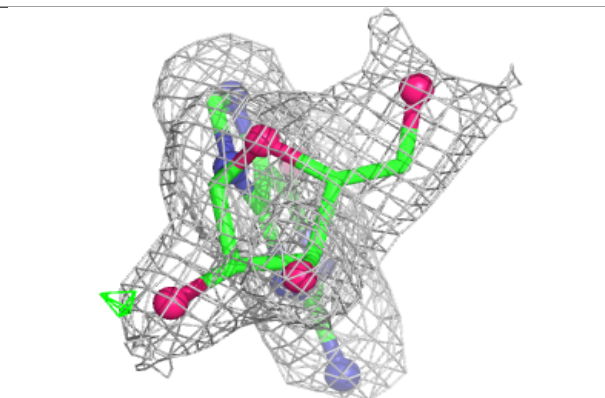
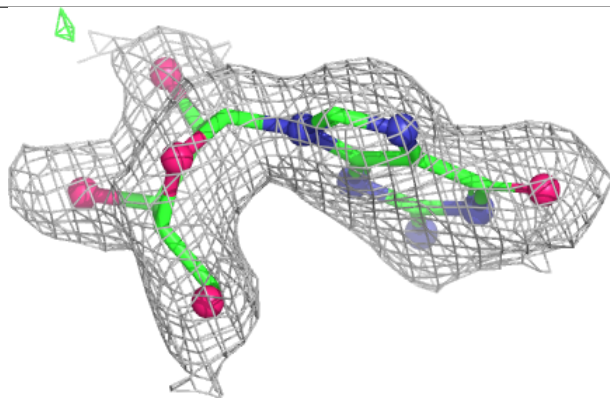
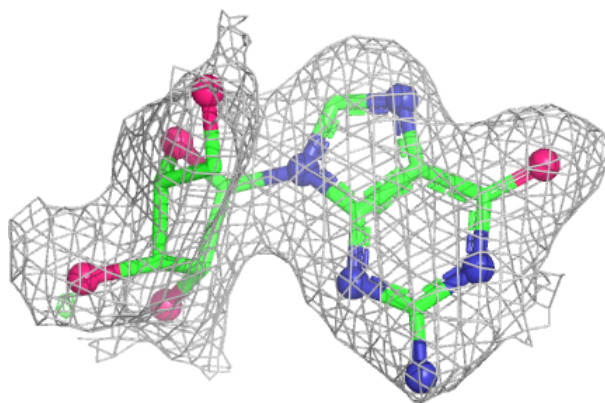
**Electron density around GMP F 1237:**

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

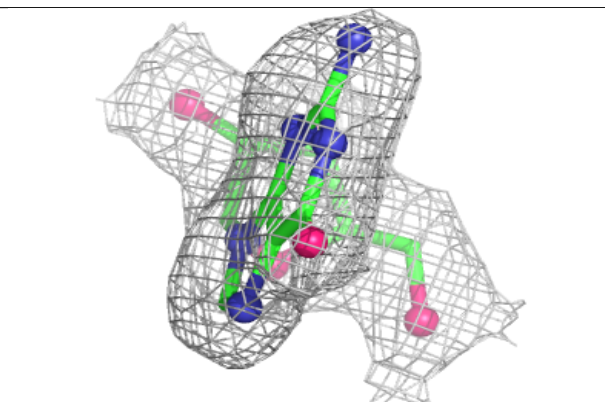
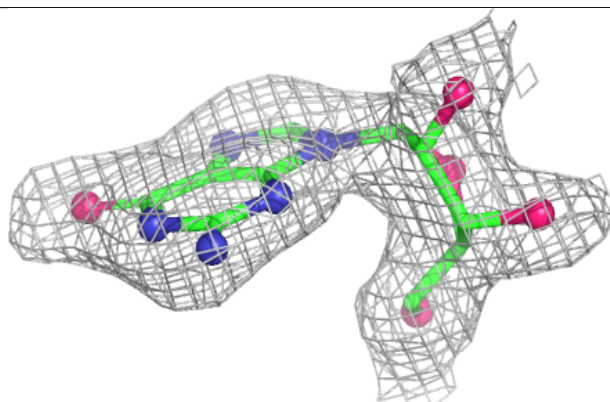
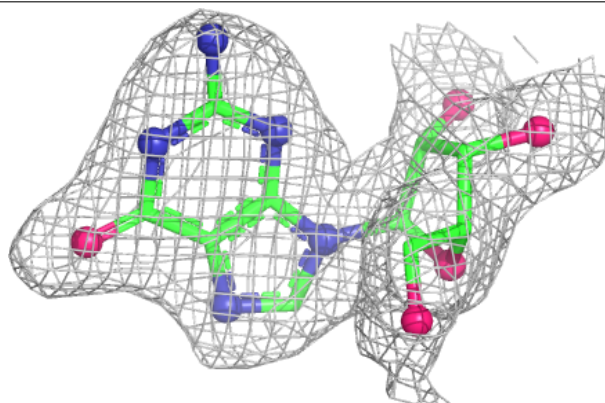


Electron density around GMP A 1237:

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

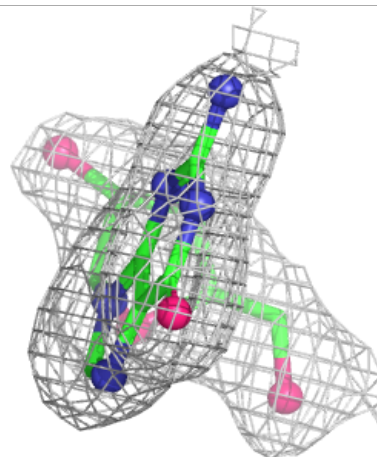
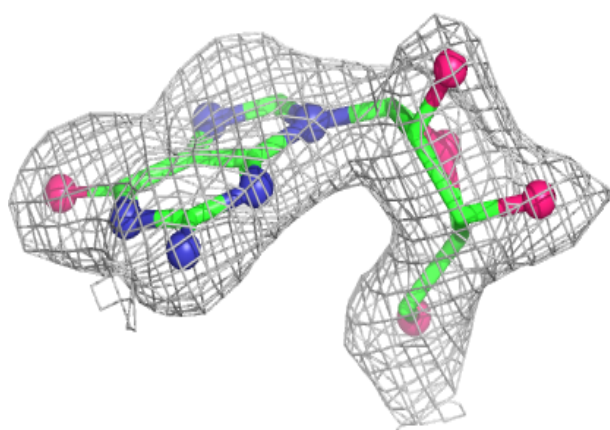
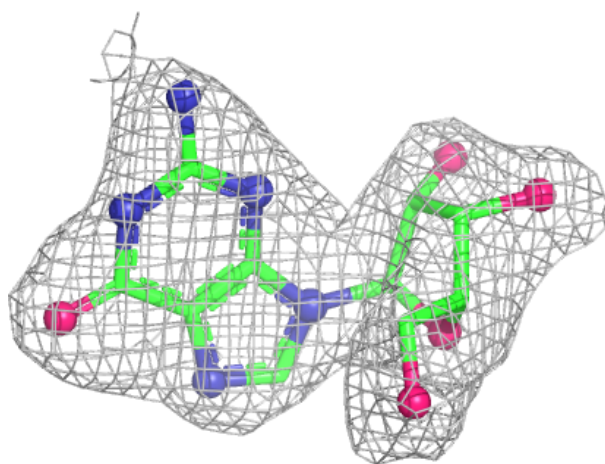
**Electron density around GMP B 1237:**

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



Electron density around GMP E 1237:

$2mF_o-DF_c$ (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.