Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 10:11 AM EDT

PDB ID	:	3DGE
Title	:	Structure of a histidine kinase-response regulator complex reveals insights into Two-component signaling and a novel cis-autophosphorylation mechanism
Authors	:	Casino, P.; Marina, A.
Deposited on	:	2008-06-13
Resolution	:	2.80 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 (i) 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 (i)) were used in the production of this report:

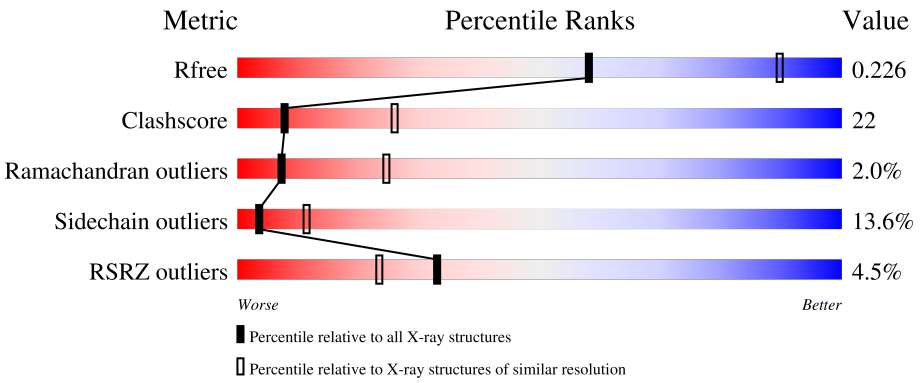
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 <=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	258	4% 45% 40% 6% • 8%
1	B	258	5% 50% 32% 9% • 8%
2	C	122	2% 56% 37% 7% •
2	D	122	5% 53% 38% 8% •

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5951 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.

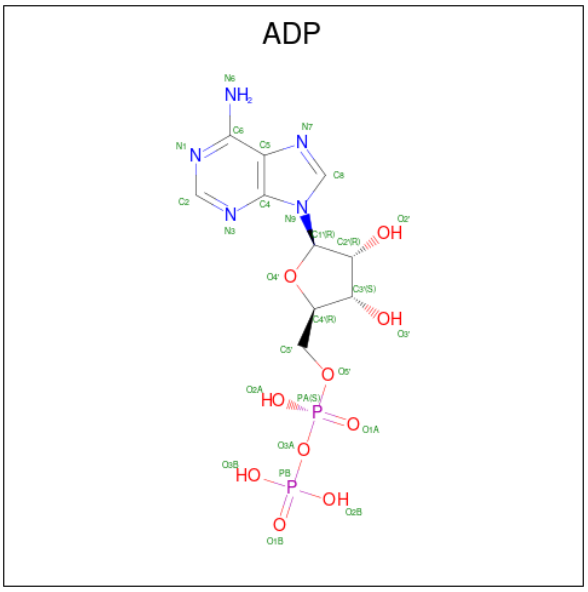
- Molecule 1 is a protein called Sensor protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	237	1886	1202	319	362	3	0	0	0
1	B	237	1886	1202	319	362	3	0	0	0

- Molecule 2 is a protein called Response regulator.

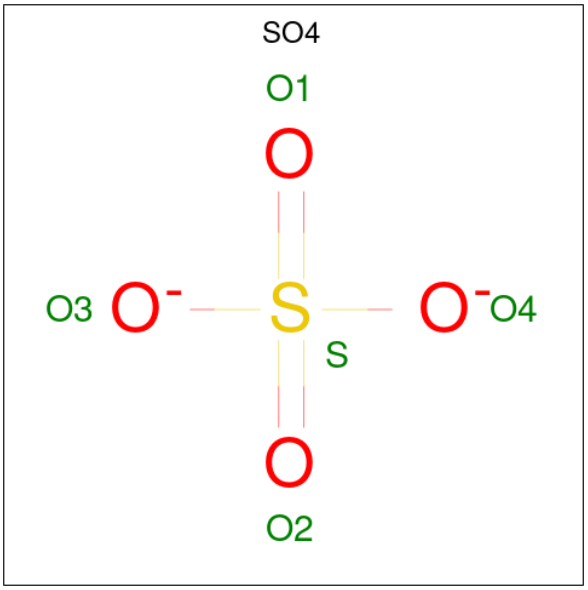
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	122	970	625	157	183	5	0	0	0
2	D	122	970	625	157	183	5	0	0	0

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



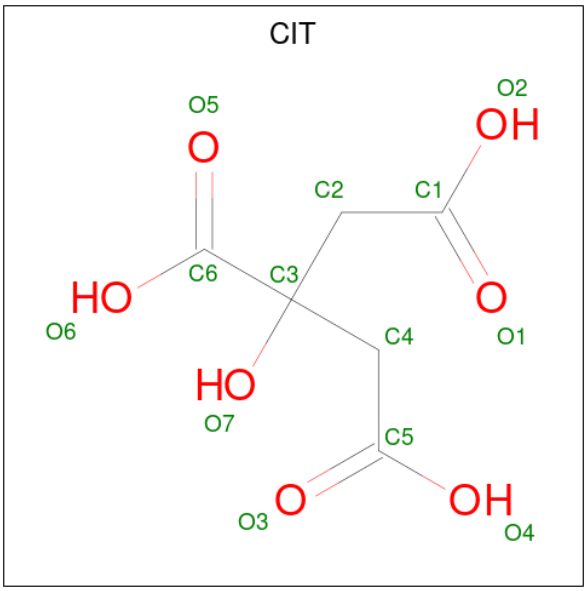
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	27	10	5	10	2	0	0
			Total	C	N	O	P		
3	B	1	27	10	5	10	2	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
4	C	1	5	4	1	0	0
			Total	O	S		
4	D	1	5	4	1	0	0

- Molecule 5 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	N	1	13	6	7	0	0

- Molecule 6 is water.

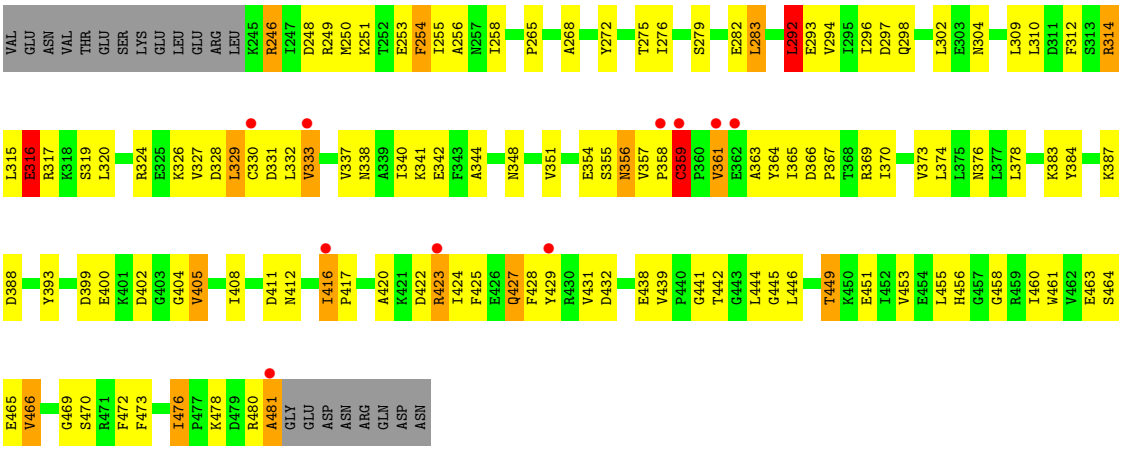
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	37	37	37	0	0
			Total	O		
6	C	27	27	27	0	0
			Total	O		
6	B	60	60	60	0	0
			Total	O		
6	D	36	36	36	0	0
			Total	O		
6	N	2	2	2	0	0

3 Residue-property plots (i)

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

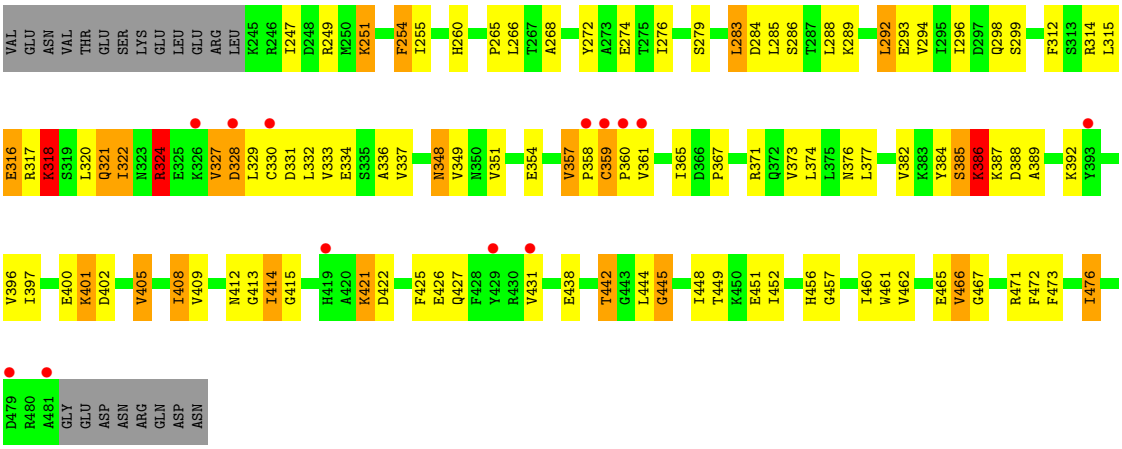
• Molecule 1: Sensor protein

Chain A:



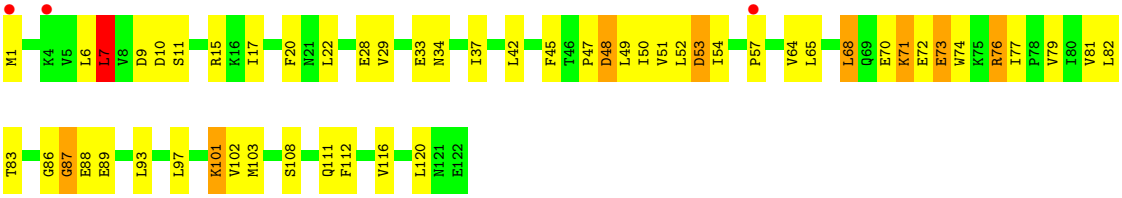
• Molecule 1: Sensor protein

Chain B:



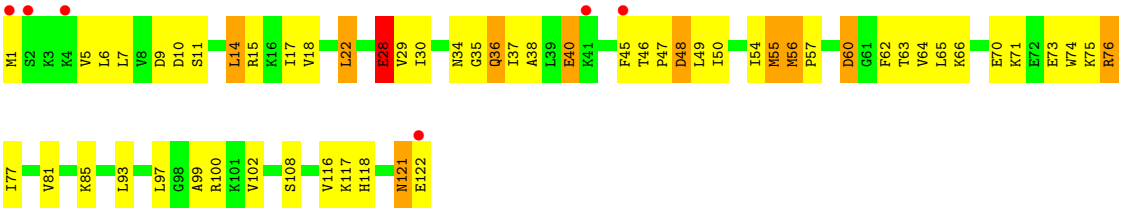
• Molecule 2: Response regulator

Chain C:



- Molecule 2: Response regulator

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	183.96Å 90.67Å 68.98Å 90.00° 108.42° 90.00°	Depositor
Resolution (Å)	43.64 – 2.80 43.63 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (43.64-2.80) 99.7 (43.63-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	1.12 (at 2.69Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.234 , 0.284 0.227 , 0.226	Depositor DCC
R_{free} test set	1502 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	67.8	Xtriage
Anisotropy	0.521	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 72.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5951	wwPDB-VP
Average B, all atoms (Å^2)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.08% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: ADP, SO4, CIT

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	1.04	7/1918 (0.4%)	0.90	6/2597 (0.2%)
1	B	1.10	14/1918 (0.7%)	0.80	2/2597 (0.1%)
2	C	0.80	4/983 (0.4%)	0.72	1/1318 (0.1%)
2	D	0.84	3/983 (0.3%)	0.81	2/1318 (0.2%)
All	All	0.99	28/5802 (0.5%)	0.82	11/7830 (0.1%)

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
1	B	0	1
All	All	0	2

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	314	ARG	CZ-NH1	26.95	1.68	1.33
1	B	354	GLU	CD-OE2	15.32	1.42	1.25
1	B	426	GLU	CD-OE1	13.76	1.40	1.25
1	B	324	ARG	CZ-NH2	12.18	1.48	1.33
1	B	354	GLU	CD-OE1	12.10	1.39	1.25
1	A	314	ARG	CZ-NH2	10.06	1.46	1.33
1	A	388	ASP	CG-OD1	8.80	1.45	1.25
2	D	76	ARG	CZ-NH1	8.72	1.44	1.33
1	B	422	ASP	CG-OD1	8.41	1.44	1.25
2	D	28	GLU	CD-OE2	8.08	1.34	1.25
1	B	465	GLU	CD-OE2	7.96	1.34	1.25
1	B	426	GLU	CD-OE2	7.63	1.34	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	316	GLU	CD-OE1	7.60	1.34	1.25
1	B	386	LYS	CE-NZ	7.31	1.67	1.49
1	B	318	LYS	CG-CD	7.27	1.77	1.52
1	A	388	ASP	CG-OD2	7.25	1.42	1.25
2	C	73	GLU	CD-OE2	7.01	1.33	1.25
2	C	89	GLU	CD-OE1	6.81	1.33	1.25
1	B	465	GLU	CD-OE1	6.76	1.33	1.25
2	D	28	GLU	CD-OE1	6.25	1.32	1.25
1	B	466	VAL	C-O	6.18	1.35	1.23
1	A	348	ASN	CG-OD1	6.08	1.37	1.24
1	A	481	ALA	C-O	5.99	1.34	1.23
2	C	73	GLU	CD-OE1	5.81	1.32	1.25
2	C	33	GLU	CD-OE2	5.64	1.31	1.25
1	B	466	VAL	C-N	5.47	1.42	1.33
1	B	422	ASP	CG-OD2	5.37	1.37	1.25
1	A	314	ARG	CD-NE	5.00	1.54	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	314	ARG	NE-CZ-NH2	-18.81	110.89	120.30
1	A	314	ARG	NE-CZ-NH1	16.38	128.49	120.30
1	B	324	ARG	NE-CZ-NH1	-12.76	113.92	120.30
1	B	324	ARG	NE-CZ-NH2	9.15	124.87	120.30
2	D	76	ARG	NE-CZ-NH2	-8.94	115.83	120.30
2	D	76	ARG	NE-CZ-NH1	7.11	123.86	120.30
1	A	388	ASP	CB-CG-OD2	-6.01	112.89	118.30
2	C	7	LEU	CA-CB-CG	5.52	128.00	115.30
1	A	314	ARG	CD-NE-CZ	-5.43	116.00	123.60
1	A	423	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	A	292	LEU	CB-CG-CD1	5.11	119.69	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	359	CYS	Peptide
1	B	359	CYS	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	1886	0	1896	107	0
1	B	1886	0	1896	93	0
2	C	970	0	1024	40	0
2	D	970	0	1024	36	0
3	A	27	0	12	4	0
3	B	27	0	12	2	0
4	C	5	0	0	1	0
4	D	5	0	0	0	0
5	N	13	0	5	0	0
6	A	37	0	0	5	0
6	B	60	0	0	3	0
6	C	27	0	0	1	0
6	D	36	0	0	2	0
6	N	2	0	0	0	0
All	All	5951	0	5869	253	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 22.

All (253) 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:B:318:LYS:CD	1:B:318:LYS:CG	1.77	1.54
1:A:314:ARG:NH1	1:A:314:ARG:CZ	1.68	1.51
1:A:314:ARG:NH2	6:A:151:HOH:O	1.83	1.12
1:A:357:VAL:HG22	1:A:358:PRO:HD2	1.31	1.11
1:A:255:ILE:HD11	1:B:254:PHE:CZ	1.97	1.00
1:A:387:LYS:HE2	2:C:57:PRO:HB3	1.43	1.00
2:C:10:ASP:OD1	6:C:177:HOH:O	1.81	0.97
1:B:376:ASN:HD21	1:B:445:GLY:C	1.70	0.94
1:A:258:ILE:HD11	1:A:309:LEU:HD21	1.49	0.94
1:A:376:ASN:ND2	1:A:449:THR:OG1	2.06	0.88
1:A:316:GLU:HG3	1:A:427:GLN:HE22	1.41	0.85
1:A:431:VAL:HB	3:A:500:ADP:O2'	1.77	0.84
1:A:255:ILE:HD11	1:B:254:PHE:HZ	1.42	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:28:GLU:OE1	2:D:28:GLU:HA	1.81	0.81
2:D:100:ARG:NH1	2:D:122:GLU:OE1	2.13	0.81
1:B:387:LYS:HE3	2:D:57:PRO:HB3	1.62	0.80
2:C:49:LEU:HD23	2:C:116:VAL:HG13	1.62	0.79
1:B:318:LYS:CG	1:B:318:LYS:CE	2.60	0.79
1:B:376:ASN:ND2	1:B:445:GLY:O	2.16	0.78
2:C:73:GLU:CD	2:C:73:GLU:H	1.87	0.78
1:B:318:LYS:CD	1:B:318:LYS:CB	2.61	0.78
2:C:101:LYS:HD3	2:C:102:VAL:N	2.00	0.76
1:A:314:ARG:CZ	6:A:151:HOH:O	2.27	0.76
1:B:365:ILE:HG22	1:B:456:HIS:HD2	1.50	0.76
1:A:255:ILE:CD1	1:B:254:PHE:HZ	1.99	0.75
1:B:414:ILE:HG22	1:B:415:GLY:H	1.51	0.74
1:B:316:GLU:HG3	1:B:427:GLN:HE22	1.51	0.74
1:B:357:VAL:HG21	1:B:400:GLU:HG3	1.68	0.74
1:A:258:ILE:CD1	1:A:309:LEU:HD21	2.20	0.72
1:B:376:ASN:ND2	1:B:449:THR:OG1	2.22	0.72
1:A:333:VAL:O	1:A:337:VAL:HG23	1.90	0.70
2:D:55:MET:O	6:D:196:HOH:O	2.08	0.70
1:B:334:GLU:HA	1:B:337:VAL:HB	1.74	0.70
1:A:327:VAL:HG11	1:A:332:LEU:HD12	1.73	0.70
1:A:428:PHE:CE2	1:B:251:LYS:HE2	2.27	0.69
1:B:337:VAL:HG13	1:B:351:VAL:CG1	2.22	0.69
1:B:327:VAL:HG11	1:B:332:LEU:HD12	1.74	0.68
1:A:315:LEU:HD21	1:A:320:LEU:HD23	1.74	0.67
1:B:324:ARG:HD3	1:B:324:ARG:H	1.57	0.67
1:A:316:GLU:OE1	1:A:427:GLN:NE2	2.28	0.67
1:B:329:LEU:O	1:B:333:VAL:HG23	1.94	0.66
1:B:387:LYS:CE	2:D:57:PRO:HB3	2.23	0.66
1:B:431:VAL:HB	3:B:600:ADP:O2'	1.94	0.66
1:A:387:LYS:CE	2:C:57:PRO:HB3	2.22	0.66
1:B:386:LYS:HE2	1:B:388:ASP:H	1.61	0.65
1:A:316:GLU:CG	1:A:427:GLN:HE22	2.08	0.65
1:A:365:ILE:HG22	1:A:456:HIS:HD2	1.62	0.65
2:D:30:ILE:HD13	2:D:45:PHE:CZ	2.32	0.64
2:D:10:ASP:OD1	6:D:196:HOH:O	2.14	0.64
2:C:49:LEU:CD2	2:C:116:VAL:HG13	2.28	0.63
1:A:428:PHE:HE2	1:B:251:LYS:HE2	1.62	0.63
2:C:82:LEU:HD23	2:C:103:MET:HB3	1.80	0.63
1:B:442:THR:HG22	1:B:444:LEU:HD13	1.80	0.62
2:D:5:VAL:O	2:D:29:VAL:HA	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:93:LEU:O	2:C:97:LEU:HB2	2.00	0.62
1:B:324:ARG:HD3	1:B:324:ARG:N	2.14	0.62
2:D:74:TRP:O	2:D:77:ILE:HG12	1.98	0.62
1:A:254:PHE:CE2	1:B:255:ILE:HD11	2.35	0.62
1:A:312:PHE:HE1	1:A:427:GLN:NE2	1.98	0.62
2:C:71:LYS:HG3	2:C:73:GLU:OE2	2.01	0.61
1:B:314:ARG:HB3	1:B:320:LEU:HD13	1.83	0.60
2:D:34:ASN:OD1	2:D:37:ILE:HG12	2.02	0.60
1:B:337:VAL:HG13	1:B:351:VAL:HG12	1.84	0.60
1:A:423:ARG:NH2	1:A:429:TYR:CE1	2.70	0.60
1:B:373:VAL:HG13	1:B:449:THR:HG23	1.84	0.59
1:A:317:ARG:HB3	1:A:319:SER:OG	2.03	0.59
2:C:49:LEU:HB2	2:C:120:LEU:HD21	1.84	0.59
1:B:321:GLN:O	1:B:322:ILE:HG13	2.03	0.58
1:B:385:SER:HA	1:B:413:GLY:HA3	1.85	0.58
1:A:272:TYR:CE1	2:C:17:ILE:HB	2.39	0.58
1:A:340:ILE:HD13	1:A:378:LEU:HB3	1.85	0.58
1:B:442:THR:CG2	1:B:444:LEU:HD13	2.34	0.58
2:C:93:LEU:O	2:C:93:LEU:HD12	2.05	0.57
1:B:330:CYS:SG	1:B:361:VAL:HG23	2.44	0.57
1:B:349:VAL:HG11	1:B:382:VAL:HG22	1.86	0.56
1:B:315:LEU:HD22	1:B:451:GLU:HG2	1.86	0.56
1:A:276:ILE:HD13	1:A:292:LEU:HD13	1.87	0.56
2:C:48:ASP:OD1	2:C:48:ASP:N	2.38	0.56
1:A:355:SER:C	1:A:356:ASN:HD22	2.09	0.56
1:B:322:ILE:HG22	1:B:324:ARG:HD2	1.88	0.56
1:A:312:PHE:CE1	1:A:427:GLN:NE2	2.74	0.56
1:A:304:ASN:HD22	1:A:304:ASN:N	2.02	0.55
1:A:316:GLU:HG3	1:A:427:GLN:NE2	2.17	0.55
1:A:328:ASP:HA	1:A:361:VAL:O	2.07	0.55
1:A:292:LEU:HD21	1:B:276:ILE:HG22	1.87	0.54
2:D:36:GLN:O	2:D:40:GLU:HG3	2.07	0.54
1:A:246:ARG:HE	1:A:250:MET:CE	2.20	0.54
1:A:365:ILE:HG22	1:A:456:HIS:CD2	2.41	0.54
1:A:329:LEU:O	1:A:333:VAL:HG22	2.07	0.54
2:C:53:ASP:OD1	4:C:150:SO4:O3	2.25	0.54
1:B:289:LYS:O	1:B:293:GLU:HG2	2.08	0.54
1:B:279:SER:O	1:B:283:LEU:HB2	2.07	0.53
1:A:420:ALA:O	1:A:424:ILE:HG23	2.09	0.53
2:C:34:ASN:OD1	2:C:37:ILE:HG12	2.08	0.53
1:B:327:VAL:HG11	1:B:332:LEU:CD1	2.37	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:337:VAL:HG13	1:B:351:VAL:HG11	1.89	0.53
2:D:63:THR:O	2:D:64:VAL:C	2.47	0.53
1:A:376:ASN:HD21	1:A:445:GLY:C	2.11	0.53
1:B:438:GLU:HA	2:D:55:MET:HG2	1.92	0.53
2:D:18:VAL:HG12	2:D:22:LEU:HD22	1.91	0.53
1:A:458:GLY:HA3	1:A:476:ILE:HD13	1.92	0.52
2:D:60:ASP:OD2	2:D:62:PHE:HB2	2.09	0.52
1:A:446:LEU:HG	3:A:500:ADP:O1A	2.09	0.52
1:A:423:ARG:NH2	1:A:429:TYR:HE1	2.08	0.52
1:B:421:LYS:NZ	6:B:153:HOH:O	2.43	0.52
1:A:340:ILE:CD1	1:A:378:LEU:HD13	2.40	0.52
1:B:333:VAL:O	1:B:337:VAL:HG23	2.10	0.52
1:A:314:ARG:CZ	6:A:152:HOH:O	2.56	0.52
1:A:429:TYR:O	3:A:500:ADP:H5'1	2.11	0.51
1:B:260:HIS:HB2	6:B:98:HOH:O	2.11	0.51
1:B:312:PHE:HB2	1:B:444:LEU:HD23	1.92	0.51
1:A:324:ARG:HD3	1:A:364:TYR:O	2.09	0.51
1:A:332:LEU:HD11	1:A:367:PRO:HA	1.91	0.51
2:C:45:PHE:CZ	2:C:47:PRO:HG3	2.45	0.51
1:A:255:ILE:CD1	1:B:254:PHE:CZ	2.76	0.51
1:A:265:PRO:HG2	1:A:302:LEU:HD13	1.93	0.51
1:B:408:ILE:HA	1:B:472:PHE:O	2.10	0.51
2:C:73:GLU:CD	2:C:73:GLU:N	2.62	0.51
2:D:93:LEU:O	2:D:97:LEU:HB2	2.11	0.51
1:A:383:LYS:HD3	1:A:439:VAL:HG13	1.93	0.51
1:A:255:ILE:HD11	1:B:254:PHE:CE2	2.43	0.50
2:C:88:GLU:CD	1:B:321:GLN:HE22	2.14	0.50
1:A:356:ASN:HD22	1:A:356:ASN:N	2.10	0.50
1:A:432:ASP:HA	6:A:125:HOH:O	2.11	0.50
2:D:49:LEU:CD2	2:D:116:VAL:HG13	2.41	0.50
2:C:108:SER:HB3	2:C:111:GLN:HB2	1.93	0.50
1:B:284:ASP:O	1:B:285:LEU:C	2.50	0.50
1:B:448:ILE:HG22	1:B:452:ILE:HD11	1.94	0.50
1:B:373:VAL:HG13	1:B:449:THR:CG2	2.41	0.50
2:C:112:PHE:O	2:C:116:VAL:HG23	2.11	0.49
1:B:389:ALA:HB3	1:B:392:LYS:HD3	1.94	0.49
1:A:384:TYR:CZ	3:A:500:ADP:H2'	2.48	0.49
2:C:101:LYS:HD3	2:C:101:LYS:C	2.29	0.49
1:B:328:ASP:HA	1:B:361:VAL:O	2.12	0.49
1:A:404:GLY:HA2	1:A:478:LYS:H	1.78	0.49
1:B:268:ALA:HB3	1:B:298:GLN:HG3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:TYR:CE1	1:A:412:ASN:HB3	2.48	0.49
2:C:9:ASP:OD1	2:C:11:SER:HB3	2.12	0.49
1:A:302:LEU:HD23	1:B:266:LEU:HD13	1.94	0.49
2:D:71:LYS:O	2:D:75:LYS:HB2	2.12	0.49
1:A:365:ILE:HD13	1:A:370:ILE:HG13	1.95	0.49
1:A:256:ALA:O	6:A:38:HOH:O	2.20	0.48
1:A:411:ASP:OD1	1:A:470:SER:N	2.43	0.48
1:B:348:ASN:N	1:B:348:ASN:HD22	2.12	0.48
2:D:71:LYS:HG3	2:D:74:TRP:CE3	2.48	0.48
2:D:6:LEU:HD23	2:D:50:ILE:HG12	1.95	0.48
1:B:333:VAL:O	1:B:337:VAL:N	2.47	0.48
1:A:315:LEU:C	1:A:317:ARG:H	2.17	0.48
1:A:480:ARG:O	1:A:481:ALA:CB	2.61	0.48
1:A:254:PHE:CZ	1:B:255:ILE:HD11	2.49	0.48
1:A:408:ILE:HA	1:A:472:PHE:O	2.13	0.48
2:D:63:THR:O	2:D:66:LYS:N	2.47	0.48
1:A:324:ARG:HH21	1:A:456:HIS:HA	1.78	0.47
1:A:464:SER:HB2	1:A:469:GLY:O	2.14	0.47
1:A:428:PHE:CD1	1:A:444:LEU:HD11	2.50	0.47
1:B:251:LYS:HA	1:B:254:PHE:HB3	1.97	0.47
1:B:316:GLU:HG3	1:B:427:GLN:NE2	2.23	0.47
1:B:401:LYS:HE3	1:B:402:ASP:H	1.80	0.47
2:D:60:ASP:OD2	2:D:62:PHE:N	2.46	0.47
2:C:54:ILE:HG12	2:C:83:THR:HB	1.97	0.47
1:B:268:ALA:CB	1:B:298:GLN:HG3	2.45	0.47
1:B:376:ASN:ND2	1:B:449:THR:HG1	2.11	0.47
1:A:366:ASP:OD2	1:A:369:ARG:HG2	2.15	0.47
2:D:14:LEU:O	2:D:18:VAL:HG23	2.14	0.47
1:A:251:LYS:HA	1:A:254:PHE:HB3	1.96	0.46
1:B:357:VAL:HG13	1:B:358:PRO:HD2	1.97	0.46
1:A:357:VAL:HG22	1:A:358:PRO:CD	2.23	0.46
1:A:480:ARG:O	1:A:481:ALA:HB3	2.16	0.46
1:A:316:GLU:CD	1:A:427:GLN:NE2	2.68	0.46
1:B:384:TYR:CZ	3:B:600:ADP:H2'	2.51	0.46
1:A:272:TYR:CZ	1:A:294:VAL:HG11	2.50	0.46
2:C:52:LEU:CD2	2:C:65:LEU:HG	2.46	0.46
1:B:272:TYR:CE1	2:D:17:ILE:HB	2.51	0.46
2:D:75:LYS:HB3	2:D:76:ARG:HD2	1.97	0.46
1:B:476:ILE:HD13	1:B:476:ILE:HA	1.75	0.46
1:B:425:PHE:CZ	1:B:460:ILE:HG12	2.50	0.45
2:D:121:ASN:C	2:D:121:ASN:HD22	2.18	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:ARG:HE	1:A:250:MET:HE3	1.80	0.45
1:A:254:PHE:HE2	1:B:255:ILE:HD11	1.79	0.45
1:A:246:ARG:HG3	1:A:249:ARG:NH2	2.31	0.45
1:A:268:ALA:HB3	1:A:298:GLN:HG3	1.97	0.45
1:A:357:VAL:CG2	1:A:358:PRO:HD2	2.22	0.45
2:C:28:GLU:HA	2:C:28:GLU:OE1	2.17	0.45
2:C:74:TRP:O	2:C:77:ILE:HG12	2.15	0.45
1:B:374:LEU:HD12	1:B:374:LEU:HA	1.85	0.45
1:B:367:PRO:O	1:B:371:ARG:HB2	2.17	0.45
2:D:30:ILE:HD13	2:D:45:PHE:HZ	1.82	0.45
1:B:396:VAL:HG22	1:B:409:VAL:HG22	1.99	0.45
2:D:48:ASP:N	2:D:48:ASP:OD1	2.50	0.45
1:A:258:ILE:HD11	1:A:309:LEU:CD2	2.33	0.45
1:A:310:LEU:HD21	1:A:314:ARG:HH21	1.82	0.45
2:C:7:LEU:HD13	2:C:15:ARG:HG2	1.99	0.45
1:B:272:TYR:CZ	1:B:294:VAL:HG11	2.52	0.45
1:A:283:LEU:HD13	2:C:20:PHE:CZ	2.52	0.45
2:C:86:GLY:O	2:C:87:GLY:C	2.55	0.45
1:B:365:ILE:HG22	1:B:456:HIS:CD2	2.40	0.45
2:D:11:SER:O	2:D:15:ARG:HG3	2.17	0.45
2:D:35:GLY:O	2:D:38:ALA:HB3	2.17	0.44
2:C:6:LEU:HB3	2:C:50:ILE:HG12	1.99	0.44
1:A:341:LYS:HA	1:A:344:ALA:HB3	1.98	0.44
1:A:373:VAL:HG13	1:A:449:THR:CG2	2.47	0.44
1:B:265:PRO:HA	1:B:298:GLN:HE21	1.83	0.44
1:A:445:GLY:O	1:A:449:THR:OG1	2.33	0.44
1:A:399:ASP:O	1:A:405:VAL:HG12	2.18	0.44
2:C:64:VAL:O	2:C:68:LEU:HD23	2.18	0.44
1:A:268:ALA:CB	1:A:298:GLN:HG3	2.48	0.43
1:A:316:GLU:CD	1:A:427:GLN:HE22	2.21	0.43
1:B:397:ILE:HB	1:B:408:ILE:HG23	2.00	0.43
1:A:326:LYS:HA	1:A:363:ALA:O	2.18	0.43
1:B:457:GLY:HA2	6:B:57:HOH:O	2.18	0.43
2:D:10:ASP:HB3	2:D:56:MET:HB3	2.00	0.43
1:A:446:LEU:HA	1:A:449:THR:OG1	2.18	0.43
2:C:7:LEU:HD23	2:C:51:VAL:HB	2.01	0.43
2:D:65:LEU:HD21	2:D:99:ALA:HB2	2.00	0.43
1:A:275:THR:CB	2:C:17:ILE:HD11	2.48	0.43
1:B:292:LEU:O	1:B:296:ILE:HG13	2.19	0.43
2:D:11:SER:HB3	2:D:14:LEU:HB2	2.01	0.43
2:C:50:ILE:HB	2:C:79:VAL:HG22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:421:LYS:O	1:B:462:VAL:CG1	2.66	0.42
1:A:310:LEU:HD21	1:A:314:ARG:NH2	2.33	0.42
1:A:416:ILE:CG2	1:A:417:PRO:HD2	2.49	0.42
1:B:386:LYS:H	1:B:413:GLY:HA2	1.84	0.42
1:B:315:LEU:C	1:B:317:ARG:H	2.21	0.42
2:D:81:VAL:HB	2:D:102:VAL:HG22	2.00	0.42
1:A:376:ASN:ND2	1:A:445:GLY:O	2.52	0.42
1:A:428:PHE:HD1	1:A:444:LEU:HD11	1.84	0.42
1:A:316:GLU:CG	1:A:427:GLN:NE2	2.80	0.42
1:A:465:GLU:HG3	1:A:466:VAL:H	1.84	0.42
1:A:315:LEU:CD2	1:A:320:LEU:HD23	2.45	0.42
1:A:449:THR:O	1:A:453:VAL:HG23	2.19	0.42
2:C:52:LEU:O	2:C:81:VAL:HA	2.20	0.42
1:B:312:PHE:CB	1:B:444:LEU:HD23	2.48	0.42
1:A:330:CYS:SG	1:A:361:VAL:HG23	2.60	0.41
1:A:425:PHE:CE2	1:A:460:ILE:HG12	2.55	0.41
2:C:71:LYS:HE3	2:C:72:GLU:H	1.85	0.41
1:B:288:LEU:HG	1:B:292:LEU:HD22	2.02	0.41
1:B:405:VAL:N	1:B:476:ILE:O	2.43	0.41
1:B:461:TRP:CE2	1:B:473:PHE:HB2	2.55	0.41
1:B:408:ILE:HD12	1:B:473:PHE:CE2	2.56	0.41
1:A:296:ILE:HD13	1:B:274:GLU:HG2	2.02	0.41
1:B:385:SER:HA	1:B:413:GLY:CA	2.50	0.41
1:A:304:ASN:N	1:A:304:ASN:ND2	2.68	0.41
2:C:9:ASP:O	2:C:15:ARG:NH1	2.46	0.41
2:C:76:ARG:HA	2:C:76:ARG:HD3	1.70	0.41
2:D:9:ASP:O	2:D:15:ARG:HD3	2.21	0.41
1:B:334:GLU:C	1:B:336:ALA:N	2.74	0.41
2:D:46:THR:HA	2:D:47:PRO:HD3	1.80	0.41
1:A:292:LEU:O	1:A:296:ILE:HG13	2.21	0.40
2:C:11:SER:O	2:C:15:ARG:HG3	2.21	0.40
1:B:473:PHE:CD1	1:B:473:PHE:N	2.89	0.40
1:A:279:SER:HB3	1:A:282:GLU:HG2	2.03	0.40
1:A:461:TRP:CE2	1:A:473:PHE:HB2	2.55	0.40
1:A:451:GLU:O	1:A:455:LEU:HG	2.21	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	235/258 (91%)	209 (89%)	21 (9%)	5 (2%)	7 23
1	B	235/258 (91%)	204 (87%)	23 (10%)	8 (3%)	3 13
2	C	120/122 (98%)	105 (88%)	14 (12%)	1 (1%)	19 49
2	D	120/122 (98%)	108 (90%)	12 (10%)	0	100 100
All	All	710/760 (93%)	626 (88%)	70 (10%)	14 (2%)	7 24

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	441	GLY
2	C	87	GLY
1	B	318	LYS
1	B	322	ILE
1	B	466	VAL
1	B	360	PRO
1	B	467	GLY
1	A	361	VAL
1	B	421	LYS
1	A	359	CYS
1	B	359	CYS
1	A	316	GLU
1	A	466	VAL
1	B	445	GLY

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	210/232 (90%)	180 (86%)	30 (14%)	3 10
1	B	210/232 (90%)	184 (88%)	26 (12%)	4 14
2	C	110/110 (100%)	98 (89%)	12 (11%)	6 19
2	D	110/110 (100%)	91 (83%)	19 (17%)	2 6
All	All	640/684 (94%)	553 (86%)	87 (14%)	3 11

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	246	ARG
1	A	248	ASP
1	A	253	GLU
1	A	254	PHE
1	A	283	LEU
1	A	292	LEU
1	A	293	GLU
1	A	297	ASP
1	A	316	GLU
1	A	329	LEU
1	A	331	ASP
1	A	333	VAL
1	A	338	ASN
1	A	342	GLU
1	A	351	VAL
1	A	354	GLU
1	A	356	ASN
1	A	359	CYS
1	A	374	LEU
1	A	400	GLU
1	A	402	ASP
1	A	405	VAL
1	A	416	ILE
1	A	422	ASP
1	A	427	GLN
1	A	438	GLU
1	A	442	THR
1	A	449	THR
1	A	463	GLU
1	A	476	ILE
2	C	1	MET
2	C	7	LEU
2	C	22	LEU

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Mol	Chain	Res	Type
2	C	29	VAL
2	C	42	LEU
2	C	48	ASP
2	C	53	ASP
2	C	68	LEU
2	C	70	GLU
2	C	71	LYS
2	C	76	ARG
2	C	101	LYS
1	B	247	ILE
1	B	249	ARG
1	B	251	LYS
1	B	254	PHE
1	B	283	LEU
1	B	286	SER
1	B	292	LEU
1	B	299	SER
1	B	321	GLN
1	B	324	ARG
1	B	327	VAL
1	B	328	ASP
1	B	331	ASP
1	B	348	ASN
1	B	357	VAL
1	B	377	LEU
1	B	385	SER
1	B	386	LYS
1	B	401	LYS
1	B	405	VAL
1	B	408	ILE
1	B	412	ASN
1	B	414	ILE
1	B	442	THR
1	B	471	ARG
1	B	476	ILE
2	D	1	MET
2	D	7	LEU
2	D	14	LEU
2	D	22	LEU
2	D	28	GLU
2	D	36	GLN
2	D	40	GLU

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Mol	Chain	Res	Type
2	D	48	ASP
2	D	54	ILE
2	D	55	MET
2	D	56	MET
2	D	60	ASP
2	D	70	GLU
2	D	73	GLU
2	D	85	LYS
2	D	108	SER
2	D	117	LYS
2	D	118	HIS
2	D	121	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	278	ASN
1	A	304	ASN
1	A	356	ASN
1	A	376	ASN
1	A	427	GLN
2	C	69	GLN
1	B	298	GLN
1	B	300	ASN
1	B	321	GLN
1	B	348	ASN
1	B	376	ASN
1	B	379	ASN
1	B	380	ASN
1	B	412	ASN
1	B	427	GLN
2	D	121	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

5 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	ADP	B	600	-	24,29,29	1.25	3 (12%)	29,45,45	1.40	3 (10%)
4	SO4	D	160	-	4,4,4	0.25	0	6,6,6	0.34	0
3	ADP	A	500	-	24,29,29	1.19	2 (8%)	29,45,45	1.56	5 (17%)
4	SO4	C	150	-	4,4,4	0.32	0	6,6,6	0.38	0
5	CIT	N	3	-	12,12,12	1.13	0	17,17,17	1.34	1 (5%)

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
5	CIT	N	3	-	-	11/16/16/16	-
3	ADP	B	600	-	-	3/12/32/32	0/3/3/3
3	ADP	A	500	-	-	4/12/32/32	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	600	ADP	PA-O3A	3.32	1.63	1.59
3	A	500	ADP	PA-O3A	2.93	1.62	1.59
3	A	500	ADP	C2-N3	2.48	1.35	1.32
3	B	600	ADP	C2-N3	2.43	1.35	1.32
3	B	600	ADP	O4'-C1'	2.40	1.44	1.40

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	500	ADP	O2A-PA-O3A	3.90	117.80	107.27
3	B	600	ADP	O2A-PA-O3A	3.76	117.44	107.27
5	N	3	CIT	O6-C6-C3	3.54	119.92	113.14
3	B	600	ADP	N3-C2-N1	-3.47	123.97	128.67
3	A	500	ADP	N3-C2-N1	-2.89	124.75	128.67
3	A	500	ADP	C4-C5-N7	-2.43	106.77	109.34
3	B	600	ADP	O3B-PB-O3A	-2.35	96.76	104.64
3	A	500	ADP	O3B-PB-O2B	2.27	116.31	107.80
3	A	500	ADP	O3B-PB-O3A	-2.03	97.81	104.64

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	500	ADP	O4'-C4'-C5'-O5'
3	B	600	ADP	PA-O3A-PB-O2B
3	B	600	ADP	O4'-C4'-C5'-O5'
5	N	3	CIT	C1-C2-C3-O7
5	N	3	CIT	C1-C2-C3-C4
5	N	3	CIT	C1-C2-C3-C6
3	A	500	ADP	C3'-C4'-C5'-O5'
3	B	600	ADP	C3'-C4'-C5'-O5'
5	N	3	CIT	C2-C3-C6-O5
5	N	3	CIT	C4-C3-C6-O6
5	N	3	CIT	O7-C3-C6-O5
5	N	3	CIT	C2-C3-C6-O6
3	A	500	ADP	PA-O3A-PB-O3B
5	N	3	CIT	O7-C3-C6-O6
5	N	3	CIT	C4-C3-C6-O5
3	A	500	ADP	PB-O3A-PA-O2A
5	N	3	CIT	O2-C1-C2-C3
5	N	3	CIT	O1-C1-C2-C3

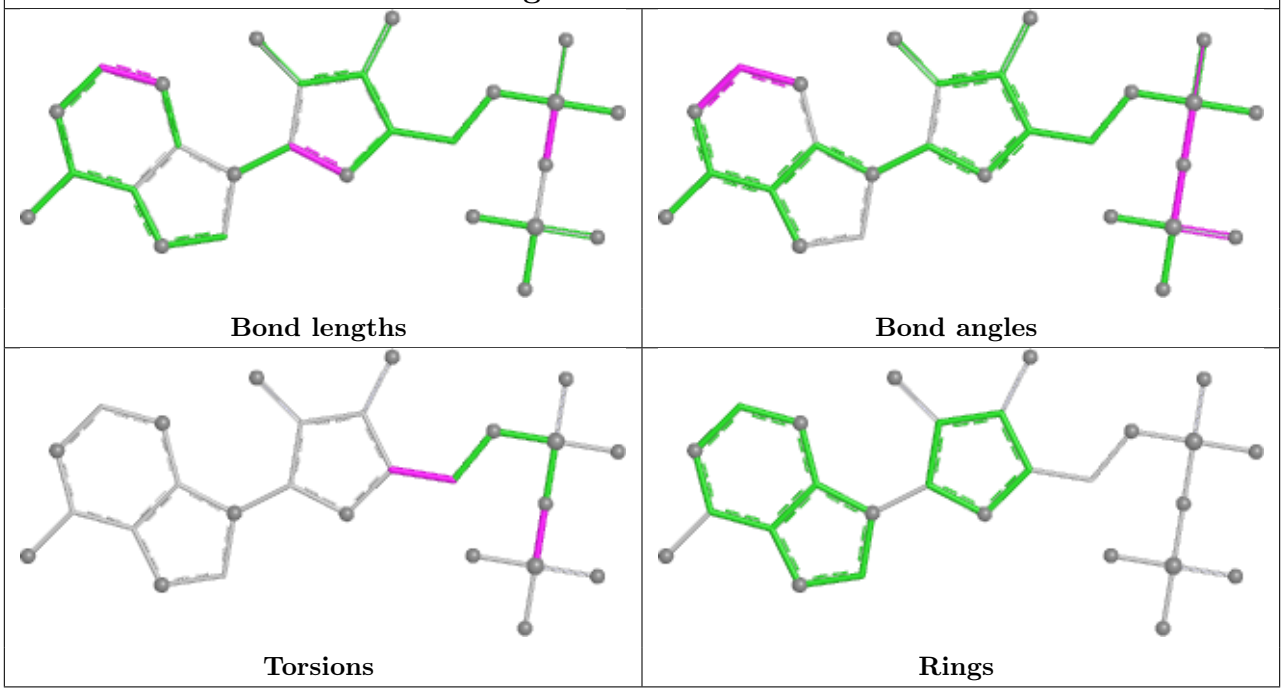
There are no ring outliers.

3 monomers are involved in 7 short contacts:

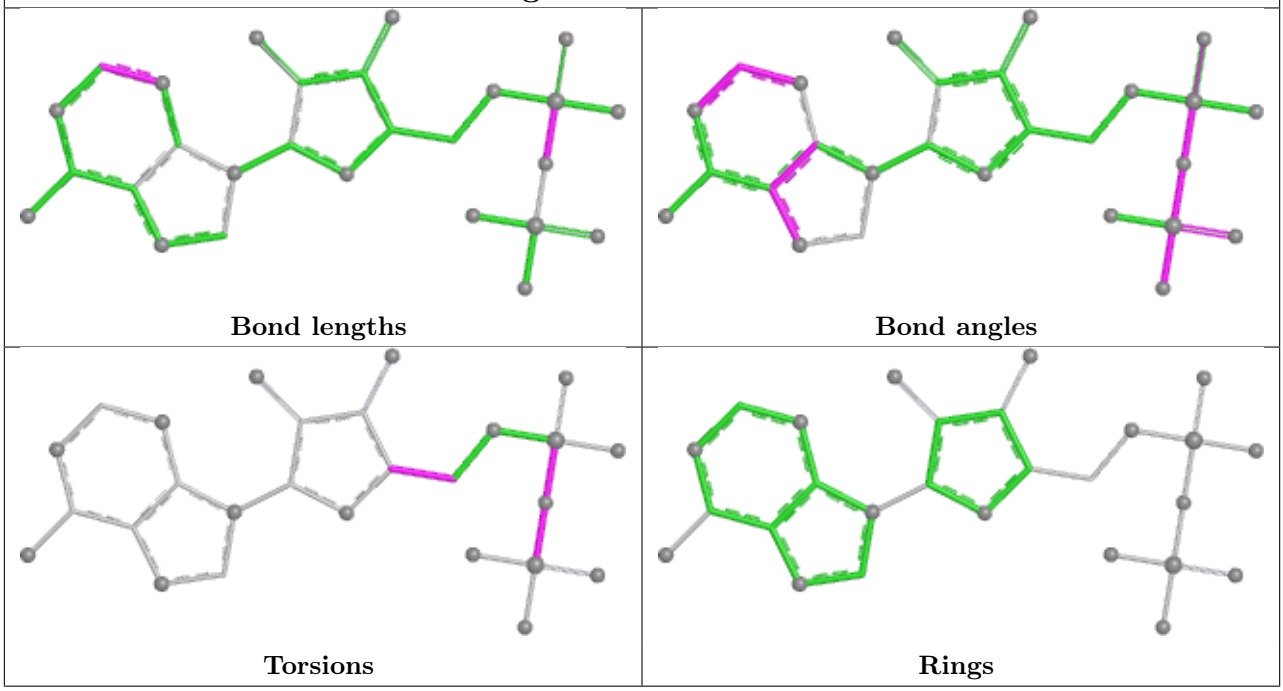
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	600	ADP	2	0
3	A	500	ADP	4	0
4	C	150	SO4	1	0

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 then 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 ADP B 600



Ligand ADP A 500



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	237/258 (91%)	0.23	10 (4%)	36	26	30, 86, 103, 108	0
1	B	237/258 (91%)	0.11	13 (5%)	25	16	33, 80, 102, 104	0
2	C	122/122 (100%)	-0.14	3 (2%)	57	47	34, 70, 86, 88	0
2	D	122/122 (100%)	0.00	6 (4%)	29	20	45, 70, 84, 87	0
All	All	718/760 (94%)	0.09	32 (4%)	33	23	30, 77, 102, 108	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	359	CYS	4.1
2	C	1	MET	4.1
1	B	361	VAL	3.9
1	B	419	HIS	3.6
1	B	393	TYR	3.2
1	B	328	ASP	3.2
1	A	330	CYS	3.1
1	B	481	ALA	3.0
2	D	1	MET	3.0
1	A	423	ARG	3.0
2	C	4	LYS	2.9
2	D	4	LYS	2.9
2	D	45	PHE	2.9
1	B	358	PRO	2.8
1	A	481	ALA	2.8
1	A	358	PRO	2.8
1	B	360	PRO	2.8
1	B	431	VAL	2.8
2	D	2	SER	2.7
1	A	362	GLU	2.7
2	D	122	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	361	VAL	2.6
1	B	330	CYS	2.5
2	C	57	PRO	2.4
2	D	41	LYS	2.4
1	A	333	VAL	2.4
1	B	359	CYS	2.3
1	B	479	ASP	2.2
1	A	416	ILE	2.1
1	B	326	LYS	2.1
1	A	429	TYR	2.0
1	B	429	TYR	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 monosaccharides in this entry.

6.4 Ligands (i)

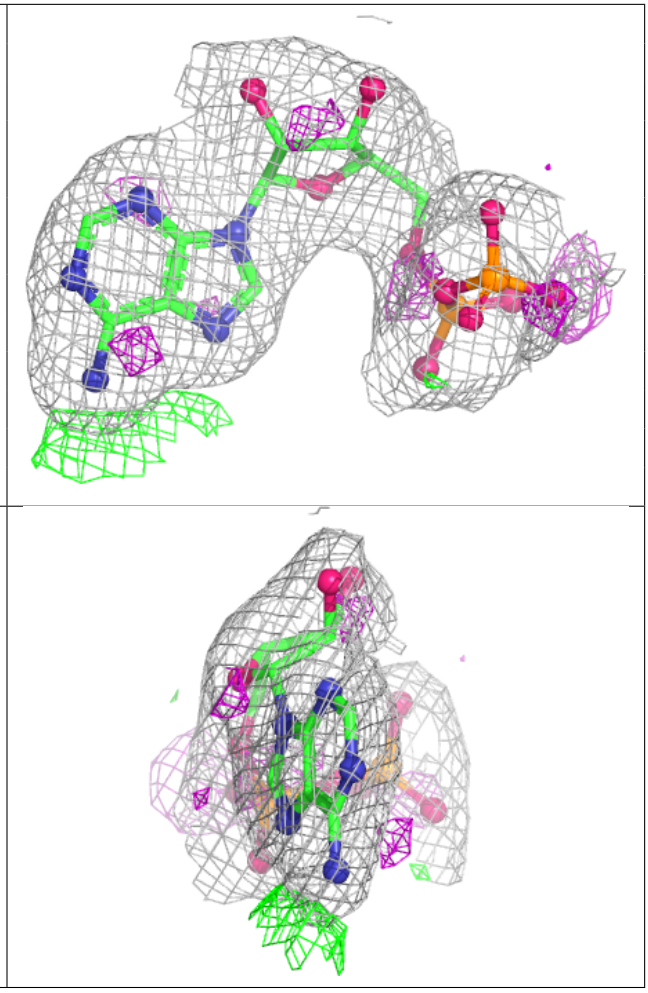
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	CIT	N	3	13/13	0.85	0.21	107,108,109,110	0
4	SO4	D	160	5/5	0.88	0.22	101,102,103,103	0
3	ADP	A	500	27/27	0.95	0.14	48,65,71,73	0
4	SO4	C	150	5/5	0.95	0.20	99,100,100,101	0
3	ADP	B	600	27/27	0.96	0.12	48,61,66,68	0

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

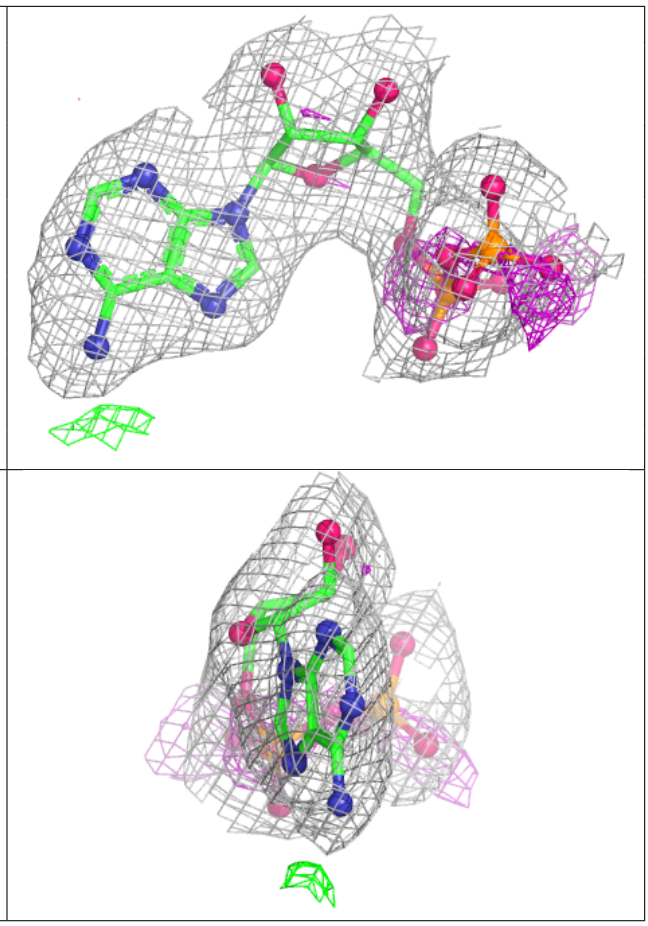
Electron density around ADP A 500:

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



Electron density around ADP B 600:

$2mF_o$-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.