Full wwPDB X-ray Structure Validation Report (i)

Aug 20, 2023 – 09:36 AM EDT

PDB ID	:	2HZY
Title	:	Mouse fumarylacetoacetate hydrolase complexes with a transition-state mimic of the complete substrate
Authors	:	Hurley, T.D.; Timm, D.E.
Deposited on	:	2006-08-09
Resolution	:	1.35 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:

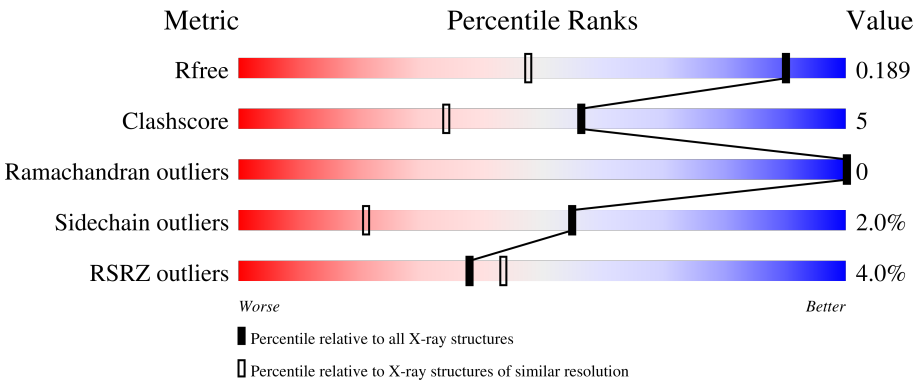
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 1.35 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	1509 (1.38-1.34)
Clashscore	141614	1551 (1.38-1.34)
Ramachandran outliers	138981	1530 (1.38-1.34)
Sidechain outliers	138945	1530 (1.38-1.34)
RSRZ outliers	127900	1487 (1.38-1.34)

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	421	5% 89% 8% • •
1	B	421	3% 89% 9% • •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	NI	A	1213	-	-	X	-

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 7532 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 Fumarylacetoacetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	416	3235	2057	562	590	26	0	4	0
1	B	418	3254	2066	570	594	24	0	5	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	cloning artifact	UNP P35505
A	0	SER	-	cloning artifact	UNP P35505
B	-1	GLY	-	cloning artifact	UNP P35505
B	0	SER	-	cloning artifact	UNP P35505

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Na		
2	A	1	1	1	0	0
2	B	1	1	1	0	0

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mn		
3	A	1	1	1	0	0
3	B	1	1	1	0	0

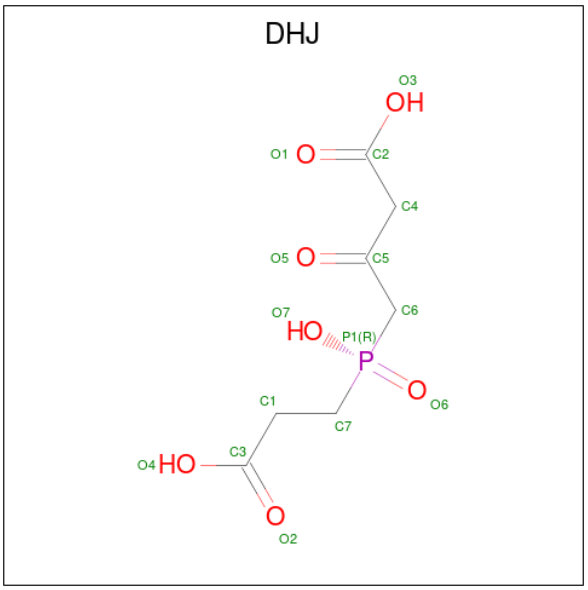
- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Ca 1 1	0	0

- Molecule 5 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	5	Total Ni 5 5	0	0
5	B	5	Total Ni 5 5	0	0

- Molecule 6 is 4-(2-CARBOXYETHYL)(HYDROXY)PHOSPHORYL]-3-OXOBUTANOIC ACID (three-letter code: DHJ) (formula: $C_7H_{11}O_7P$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O P 15 7 7 1	0	0
6	B	1	Total C O P 15 7 7 1	0	0

- Molecule 7 is water.

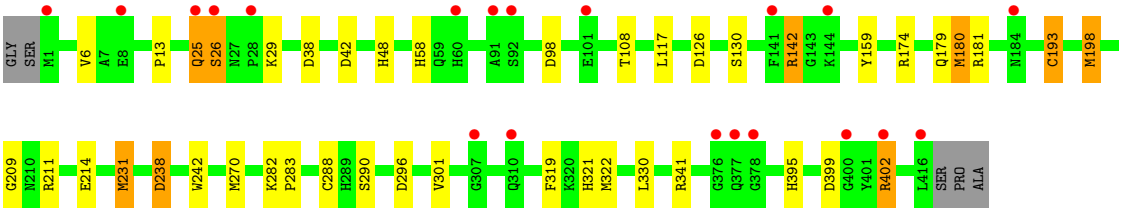
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	467	Total O 467 467	0	0
7	B	529	Total O 531 531	0	2

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.

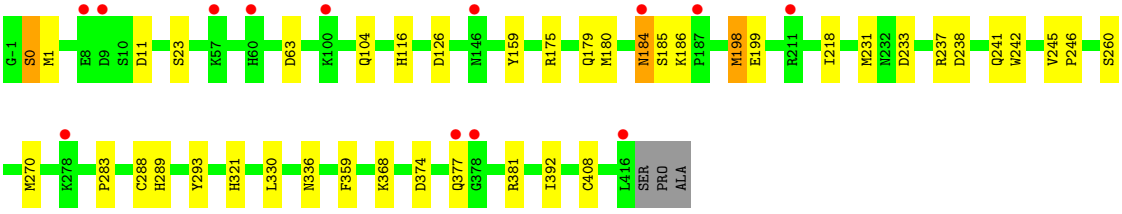
- Molecule 1: Fumarylacetoacetase

Chain A:



- Molecule 1: Fumarylacetoacetase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	64.14Å 109.47Å 67.49Å 90.00° 102.35° 90.00°	Depositor
Resolution (Å)	24.14 – 1.35 24.14 – 1.35	Depositor EDS
% Data completeness (in resolution range)	94.6 (24.14-1.35) 94.6 (24.14-1.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	2.74 (at 1.35Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.168 , 0.188 0.169 , 0.189	Depositor DCC
R_{free} test set	9466 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	12.4	Xtriage
Anisotropy	0.703	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 50.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.026 for l,-k,h	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	7532	wwPDB-VP
Average B, all atoms (Å^2)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% 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: MN, CA, NI, NA, DHJ

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	0.84	4/3343 (0.1%)	0.96	11/4537 (0.2%)
1	B	0.86	4/3366 (0.1%)	0.94	8/4570 (0.2%)
All	All	0.85	8/6709 (0.1%)	0.95	19/9107 (0.2%)

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 (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	198	MET	SD-CE	-12.40	1.08	1.77
1	A	231	MET	SD-CE	-10.73	1.17	1.77
1	B	231	MET	SD-CE	-6.73	1.40	1.77
1	A	193[A]	CYS	CB-SG	-6.09	1.72	1.82
1	A	193[B]	CYS	CB-SG	-6.09	1.72	1.82
1	B	260	SER	CB-OG	5.57	1.49	1.42
1	B	0	SER	C-N	5.13	1.45	1.34
1	A	142	ARG	CB-CG	5.01	1.66	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	231	MET	CG-SD-CE	10.25	116.59	100.20

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	238	ASP	CB-CG-OD1	7.22	124.80	118.30
1	B	238	ASP	CB-CG-OD1	6.75	124.38	118.30
1	B	374	ASP	CB-CG-OD2	6.55	124.20	118.30
1	B	11	ASP	CB-CG-OD1	6.53	124.17	118.30
1	B	175	ARG	NE-CZ-NH1	6.51	123.56	120.30
1	A	38	ASP	CB-CG-OD2	6.43	124.08	118.30
1	B	175	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	A	98	ASP	CB-CG-OD2	5.89	123.60	118.30
1	A	296	ASP	CB-CG-OD1	5.78	123.50	118.30
1	A	142	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	42	ASP	CB-CG-OD1	5.33	123.10	118.30
1	A	181	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	A	174	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	231	MET	CG-SD-CE	-5.26	91.78	100.20
1	B	381	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	B	359	PHE	CB-CG-CD2	5.13	124.39	120.80
1	A	341	ARG	NE-CZ-NH2	-5.12	117.74	120.30
1	A	319	PHE	CB-CG-CD1	5.01	124.31	120.80

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	209	GLY	Peptide
1	B	0	SER	Mainchain

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	3235	0	3183	37	0
1	B	3254	0	3193	25	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	5	0	0	3	0
5	B	5	0	0	0	0
6	A	15	0	8	0	0
6	B	15	0	8	0	0
7	A	467	0	0	15	0
7	B	531	0	0	7	0
All	All	7532	0	6392	65	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 5.

All (65) 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:231:MET:CE	1:A:231:MET:SD	1.17	1.26
1:A:282:LYS:HE3	7:A:1549:HOH:O	1.38	1.20
1:B:198:MET:CE	1:B:198:MET:SD	1.08	1.18
1:A:231:MET:SD	1:A:231:MET:HE3	1.75	1.15
1:A:231:MET:SD	1:A:231:MET:HE2	1.75	1.10
1:A:231:MET:SD	1:A:231:MET:HE1	1.75	1.09
1:B:198:MET:SD	1:B:198:MET:HE2	1.67	1.08
1:B:198:MET:SD	1:B:198:MET:HE3	1.67	1.08
1:B:198:MET:CE	1:B:198:MET:CG	2.33	1.05
1:B:198:MET:SD	1:B:198:MET:HE1	1.67	1.05
1:A:321:HIS:CE1	7:A:1520:HOH:O	2.11	1.03
1:B:321:HIS:NE2	7:B:1729:HOH:O	2.10	0.80
1:A:321:HIS:NE2	7:A:1520:HOH:O	2.11	0.80
1:A:231:MET:CE	1:A:231:MET:CG	2.60	0.79
1:A:130:SER:HB2	1:A:322[B]:MET:HE1	1.64	0.78
1:A:130:SER:HB2	1:A:322[B]:MET:CE	2.13	0.77
1:A:58:HIS:HE1	7:A:1646:HOH:O	1.68	0.76
1:A:142:ARG:HD2	7:A:1253:HOH:O	1.86	0.75
1:A:179:GLN:HE21	1:A:242:TRP:HE1	1.36	0.72
5:A:1213:NI:NI	7:A:1591:HOH:O	1.32	0.72
1:B:179:GLN:HE21	1:B:242:TRP:HE1	1.39	0.71
1:B:283:PRO:HG2	1:B:288[B]:CYS:SG	2.30	0.71
5:A:1212:NI:NI	7:A:1668:HOH:O	1.34	0.70
1:A:399:ASP:OD2	7:A:1668:HOH:O	2.10	0.69
1:B:321:HIS:CE1	7:B:1729:HOH:O	2.43	0.69
5:A:1213:NI:NI	7:A:1634:HOH:O	1.37	0.68
1:A:282:LYS:NZ	1:A:288:CYS:HB2	2.08	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:HG2	7:B:1616:HOH:O	1.94	0.67
1:A:25:GLN:HG3	1:A:26:SER:N	2.12	0.65
1:A:395:HIS:CD2	1:A:402:ARG:HD3	2.32	0.64
1:A:282:LYS:HD3	1:A:283:PRO:O	1.98	0.64
1:A:58:HIS:CE1	7:A:1646:HOH:O	2.47	0.62
1:A:179:GLN:NE2	1:A:242:TRP:HE1	1.98	0.61
1:B:179:GLN:NE2	1:B:242:TRP:HE1	1.98	0.60
1:A:270:MET:HA	1:A:270:MET:HE3	1.84	0.59
1:A:117:LEU:HD11	7:A:1329:HOH:O	2.06	0.56
1:A:108:THR:HG22	7:A:1543:HOH:O	2.06	0.55
1:B:23:SER:OG	1:B:116:HIS:HE1	1.91	0.54
1:A:282:LYS:HZ2	1:A:288:CYS:HB2	1.72	0.54
1:A:130:SER:CB	1:A:322[B]:MET:HE1	2.37	0.53
1:A:282:LYS:HZ1	1:A:288:CYS:HB2	1.75	0.52
1:A:179:GLN:O	1:A:180[B]:MET:HG3	2.10	0.52
1:B:241[A]:GLN:HE22	1:B:368:LYS:NZ	2.08	0.51
1:A:198[B]:MET:SD	1:A:301:VAL:HG11	2.51	0.50
1:B:116:HIS:CD2	7:B:1653:HOH:O	2.64	0.49
1:A:6:VAL:HG11	1:A:13:PRO:HA	1.94	0.49
1:A:48:HIS:CE1	7:A:1591:HOH:O	2.65	0.48
1:B:184:ASN:HD22	1:B:185:SER:N	2.12	0.47
1:A:126:ASP:HB3	1:A:159:TYR:CE2	2.51	0.46
1:B:186:LYS:HD2	7:B:1662:HOH:O	2.17	0.45
1:B:270:MET:HE3	1:B:270:MET:HA	2.00	0.44
1:A:282:LYS:HZ1	1:A:288:CYS:CB	2.30	0.43
7:A:1592:HOH:O	1:B:336:ASN:HB3	2.18	0.43
1:B:199:GLU:HB2	1:B:233:ASP:HB3	2.00	0.42
1:B:245:VAL:HA	1:B:246:PRO:HA	1.88	0.42
1:A:290:SER:HB2	7:A:1657:HOH:O	2.19	0.42
1:B:104:GLN:NE2	7:B:1591:HOH:O	2.52	0.42
1:B:237:ARG:O	1:B:241[A]:GLN:HG3	2.20	0.42
1:A:180[B]:MET:HE3	1:A:238:ASP:OD2	2.20	0.42
1:A:117:LEU:HD12	7:B:1477:HOH:O	2.20	0.42
1:B:218:ILE:HD11	1:B:289[A]:HIS:CD2	2.56	0.41
1:A:211:ARG:HE	1:A:214:GLU:CD	2.24	0.40
1:B:126:ASP:HB3	1:B:159:TYR:CE2	2.56	0.40
1:B:392:ILE:HB	1:B:408:CYS:HB3	2.04	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	418/421 (99%)	410 (98%)	8 (2%)	0	100	100
1	B	421/421 (100%)	408 (97%)	13 (3%)	0	100	100
All	All	839/842 (100%)	818 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	355/354 (100%)	344 (97%)	11 (3%)	40	8
1	B	357/354 (101%)	351 (98%)	6 (2%)	60	28
All	All	712/708 (101%)	695 (98%)	17 (2%)	55	15

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	26	SER
1	A	29	LYS
1	A	180[A]	MET
1	A	180[B]	MET
1	A	193[A]	CYS
1	A	193[B]	CYS
1	A	198[A]	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	198[B]	MET
1	A	330	LEU
1	A	402	ARG
1	B	63	ASP
1	B	180	MET
1	B	184	ASN
1	B	293	TYR
1	B	330	LEU
1	B	377	GLN

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

Mol	Chain	Res	Type
1	A	39	GLN
1	A	179	GLN
1	A	321	HIS
1	A	397	GLN
1	B	116	HIS
1	B	179	GLN
1	B	184	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)

Of 17 ligands modelled in this entry, 15 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
6	DHJ	B	1101	3	11,14,14	2.56	3 (27%)	11,19,19	1.43	1 (9%)
6	DHJ	A	1102	3	11,14,14	2.30	3 (27%)	11,19,19	1.87	2 (18%)

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
6	DHJ	B	1101	3	-	6/14/15/15	-
6	DHJ	A	1102	3	-	7/14/15/15	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	1101	DHJ	C7-C1	-6.93	1.32	1.52
6	A	1102	DHJ	C7-C1	-6.13	1.34	1.52
6	B	1101	DHJ	O4-C3	-2.84	1.21	1.30
6	A	1102	DHJ	O2-C3	2.36	1.30	1.22
6	A	1102	DHJ	O4-C3	-2.28	1.23	1.30
6	B	1101	DHJ	O5-C5	2.14	1.25	1.21

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	A	1102	DHJ	C7-C1-C3	4.92	121.11	113.49
6	B	1101	DHJ	C7-C1-C3	3.82	119.41	113.49
6	A	1102	DHJ	C6-P1-C7	-2.65	103.11	107.71

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1102	DHJ	C2-C4-C5-O5
6	B	1101	DHJ	C2-C4-C5-C6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
6	B	1101	DHJ	C2-C4-C5-O5
6	A	1102	DHJ	C1-C7-P1-O7
6	A	1102	DHJ	O3-C2-C4-C5
6	B	1101	DHJ	O3-C2-C4-C5
6	A	1102	DHJ	C1-C7-P1-O6
6	B	1101	DHJ	C5-C6-P1-O6
6	A	1102	DHJ	C1-C7-P1-C6
6	A	1102	DHJ	C2-C4-C5-C6
6	A	1102	DHJ	O1-C2-C4-C5
6	B	1101	DHJ	O1-C2-C4-C5
6	B	1101	DHJ	C5-C6-P1-C7

There are no ring outliers.

No monomer is involved in short contacts.

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	416/421 (98%)	0.25	20 (4%)	30	35	8, 15, 29, 39	1 (0%)
1	B	418/421 (99%)	0.07	13 (3%)	49	56	9, 14, 26, 34	0
All	All	834/842 (99%)	0.16	33 (3%)	38	43	8, 15, 28, 39	1 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	307	GLY	7.4
1	A	141	PHE	5.2
1	B	378	GLY	4.8
1	A	416	LEU	4.6
1	A	28	PRO	4.5
1	B	416	LEU	4.1
1	A	1	MET	3.8
1	A	144	LYS	3.4
1	A	400	GLY	3.3
1	A	25	GLN	3.3
1	A	378	GLY	3.2
1	B	184	ASN	3.2
1	A	60	HIS	2.9
1	B	57	LYS	2.8
1	A	91	ALA	2.6
1	A	377	GLN	2.5
1	B	187	PRO	2.5
1	B	211	ARG	2.4
1	A	92	SER	2.4
1	A	310	GLN	2.4
1	B	8	GLU	2.4
1	B	278	LYS	2.4
1	A	184	ASN	2.3
1	B	60	HIS	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	9	ASP	2.3
1	A	402	ARG	2.3
1	B	146	ASN	2.3
1	A	101	GLU	2.3
1	A	26	SER	2.3
1	A	8	GLU	2.2
1	A	376	GLY	2.2
1	B	100	LYS	2.1
1	B	377	GLN	2.1

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

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

6.3 Carbohydrates (i)

There are no 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	NI	B	1215	1/1	0.89	0.05	28,28,28,28	1
5	NI	A	1214	1/1	0.94	0.05	24,24,24,24	1
5	NI	A	1211	1/1	0.96	0.05	23,23,23,23	1
5	NI	A	1213	1/1	0.97	0.07	22,22,22,22	1
5	NI	B	1210	1/1	0.98	0.04	16,16,16,16	1
5	NI	B	1205	1/1	0.98	0.06	16,16,16,16	0
6	DHJ	A	1102	15/15	0.98	0.06	9,12,19,22	0
6	DHJ	B	1101	15/15	0.98	0.08	10,11,18,19	0
5	NI	A	1212	1/1	0.99	0.04	17,17,17,17	1
5	NI	B	1207	1/1	0.99	0.04	16,16,16,16	1
4	CA	A	1206	1/1	1.00	0.05	10,10,10,10	0
5	NI	A	1209	1/1	1.00	0.03	15,15,15,15	1
5	NI	B	1208	1/1	1.00	0.05	12,12,12,12	1
2	NA	A	1201	1/1	1.00	0.07	9,9,9,9	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	NA	B	1202	1/1	1.00	0.06	9,9,9,9	0
3	MN	A	1203	1/1	1.00	0.05	9,9,9,9	0
3	MN	B	1204	1/1	1.00	0.06	8,8,8,8	0

6.5 Other polymers (i)

There are no such residues in this entry.