Full wwPDB X-ray Structure Validation Report (i)

Sep 19, 2023 – 10:55 AM EDT

PDB ID	:	5HN8
Title	:	Crystal structure of Plasmodium vivax geranylgeranylpyrophosphate synthase complexed with BPH-1182
Authors	:	Liu, Y.-L.; Zhang, Y.; Oldfield, E.
Deposited on	:	2016-01-18
Resolution	:	2.70 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:

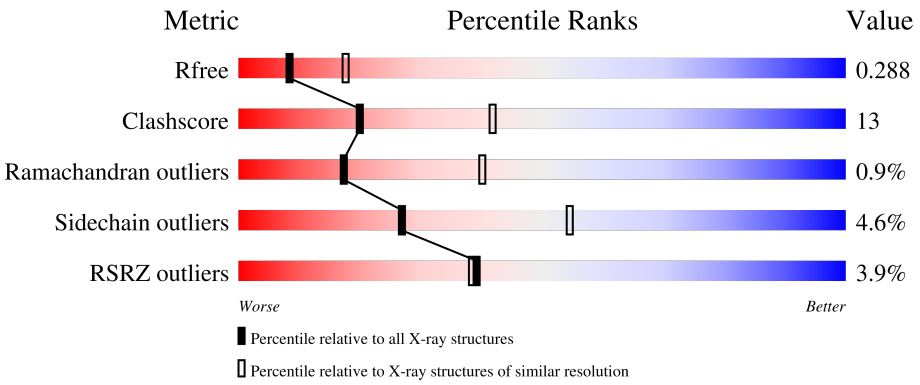
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	375	4% 66% 18% • 14%
1	B	375	3% 69% 16% • 13%
1	C	375	5% 65% 21% • 11%
1	D	375	2% 68% 19% • 11%

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
2	HXK	D	401	-	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 10771 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.

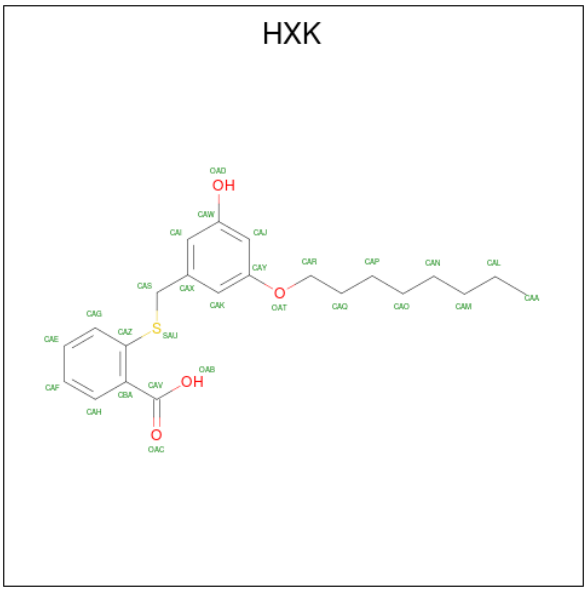
- Molecule 1 is a protein called Farnesyl pyrophosphate synthase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	321	2580	1682	405	479	14	0	0	0
1	B	326	2595	1684	414	483	14	0	0	0
1	C	334	2662	1730	426	491	15	0	0	0
1	D	334	2660	1731	424	490	15	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

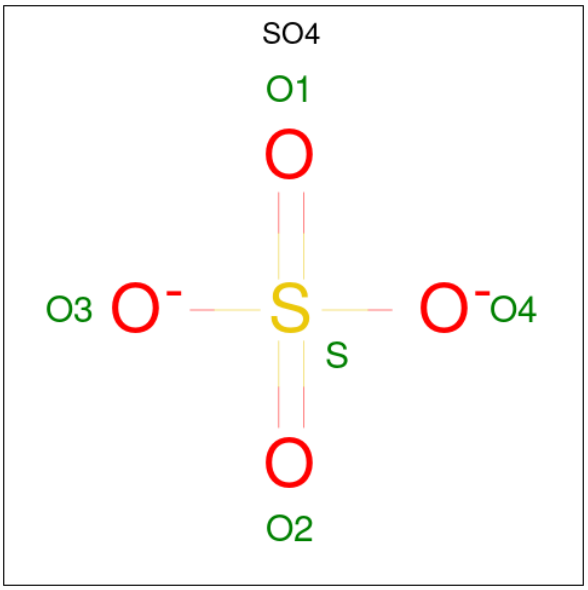
Chain	Residue	Modelled	Actual	Comment	Reference
A	134	MET	THR	SEE REMARK 999	UNP A5K4U6
A	227	ASP	ASN	SEE REMARK 999	UNP A5K4U6
B	134	MET	THR	SEE REMARK 999	UNP A5K4U6
B	227	ASP	ASN	SEE REMARK 999	UNP A5K4U6
C	134	MET	THR	SEE REMARK 999	UNP A5K4U6
C	227	ASP	ASN	SEE REMARK 999	UNP A5K4U6
D	134	MET	THR	SEE REMARK 999	UNP A5K4U6
D	227	ASP	ASN	SEE REMARK 999	UNP A5K4U6

- Molecule 2 is 2-{[3-hydroxy-5-(octyloxy)benzyl]sulfanyl}benzoic acid (three-letter code: HXK) (formula: $C_{22}H_{28}O_4S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	S		
2	A	1	27	22	4	1	0	0
2	B	1	27	22	4	1	0	0
2	C	1	27	22	4	1	0	0
2	D	1	27	22	4	1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 4 is water.

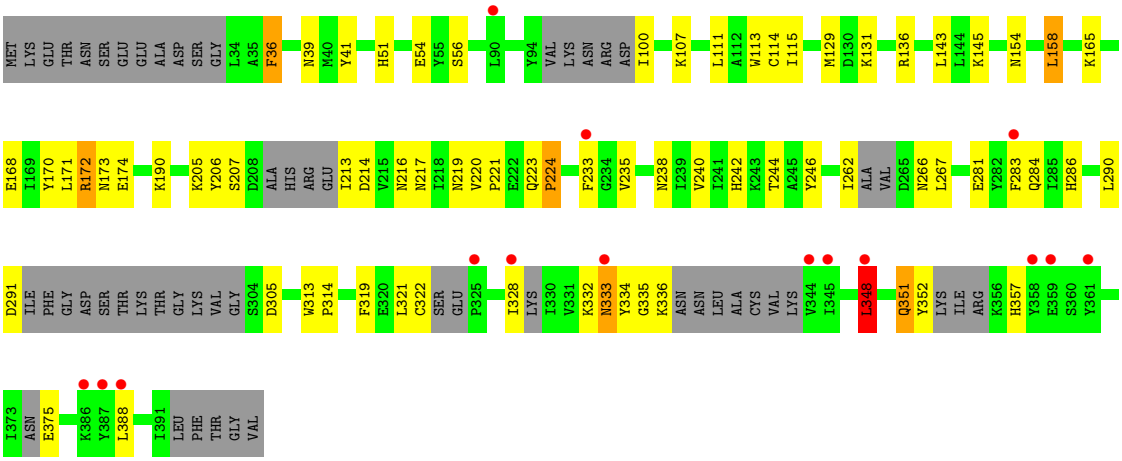
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	60	60	60	0	0
			Total	O		
4	B	19	19	19	0	0
			Total	O		
4	C	38	38	38	0	0
			Total	O		
4	D	44	44	44	0	0

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.

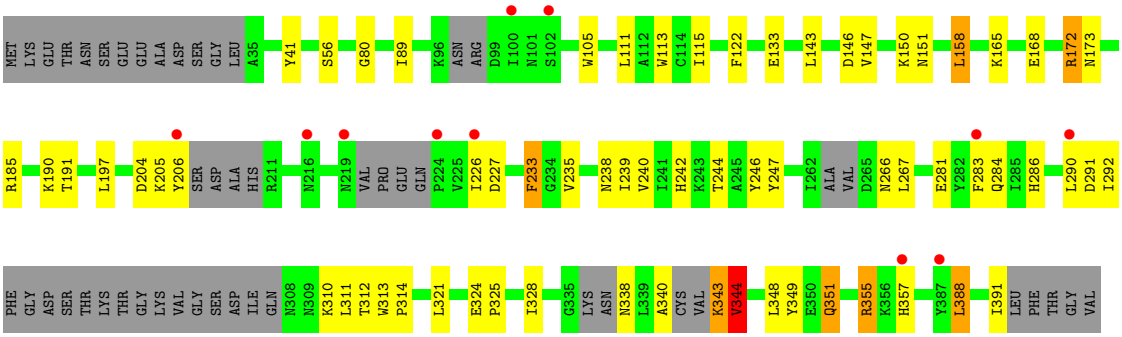
• Molecule 1: Farnesyl pyrophosphate synthase, putative

Chain A:



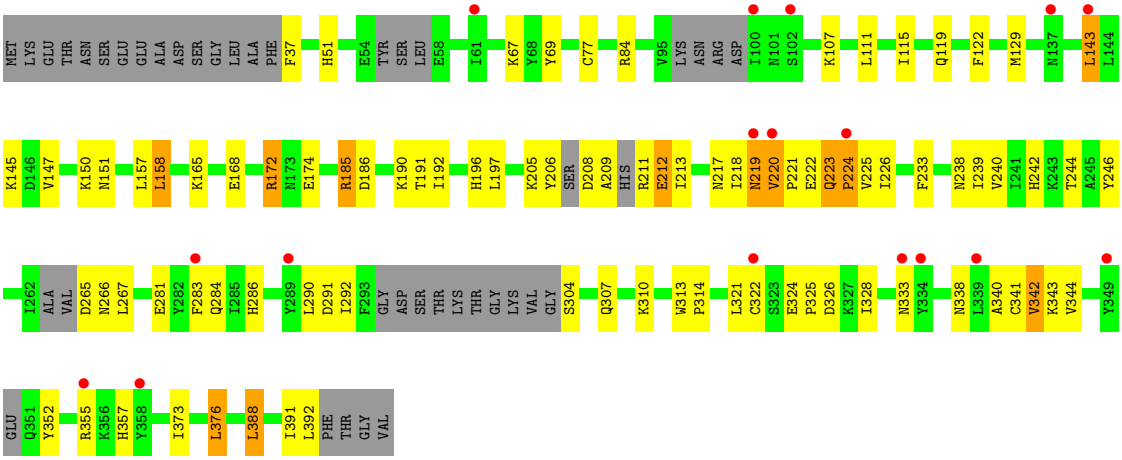
• Molecule 1: Farnesyl pyrophosphate synthase, putative

Chain B:



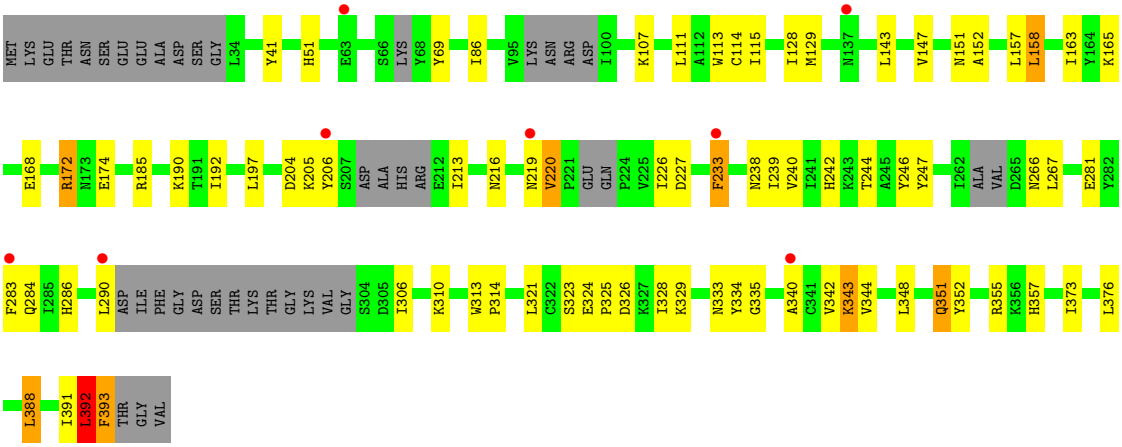
• Molecule 1: Farnesyl pyrophosphate synthase, putative

Chain C:



- Molecule 1: Farnesyl pyrophosphate synthase, putative

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	84.07Å 117.22Å 92.79Å 90.00° 116.20° 90.00°	Depositor
Resolution (Å)	31.10 – 2.70 31.10 – 2.70	Depositor EDS
% Data completeness (in resolution range)	92.0 (31.10-2.70) 91.9 (31.10-2.70)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.92 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.259 , 0.296 0.259 , 0.288	Depositor DCC
R_{free} test set	2075 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å^2)	46.1	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 54.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	10771	wwPDB-VP
Average B, all atoms (Å^2)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.87% 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: SO4, HXK

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	0.69	1/2631 (0.0%)	0.69	3/3559 (0.1%)
1	B	0.61	0/2645	0.66	1/3582 (0.0%)
1	C	0.63	1/2712 (0.0%)	0.65	2/3673 (0.1%)
1	D	0.60	0/2712	0.67	1/3675 (0.0%)
All	All	0.63	2/10700 (0.0%)	0.67	7/14489 (0.0%)

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	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	77	CYS	CB-SG	-6.57	1.71	1.82
1	A	114	CYS	CB-SG	-5.39	1.73	1.81

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	376	LEU	CA-CB-CG	5.94	128.97	115.30
1	B	344	VAL	N-CA-C	-5.78	95.38	111.00
1	C	158	LEU	CA-CB-CG	-5.76	102.05	115.30
1	A	158	LEU	CA-CB-CG	-5.74	102.10	115.30
1	A	348	LEU	CA-CB-CG	5.42	127.77	115.30
1	D	158	LEU	CA-CB-CG	-5.38	102.93	115.30
1	A	351	GLN	CB-CA-C	-5.20	100.01	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	343	LYS	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	2580	0	2433	60	0
1	B	2595	0	2431	56	0
1	C	2662	0	2528	82	0
1	D	2660	0	2523	82	0
2	A	27	0	0	3	0
2	B	27	0	0	6	0
2	C	27	0	0	2	0
2	D	27	0	0	13	0
3	D	5	0	0	0	0
4	A	60	0	0	11	0
4	B	19	0	0	4	0
4	C	38	0	0	4	0
4	D	44	0	0	5	0
All	All	10771	0	9915	258	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 13.

All (258) 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)
2:A:401:HXK:CAA	1:C:157:LEU:HD12	1.85	1.06
1:C:283:PHE:CE1	1:C:392:LEU:HG	1.90	1.05
1:D:283:PHE:CD2	2:D:401:HXK:CAH	2.41	1.02
1:A:335:GLY:HA3	4:A:501:HOH:O	1.61	1.01
1:A:291:ASP:OD2	1:A:305:ASP:HB2	1.67	0.95
1:D:283:PHE:CE2	2:D:401:HXK:CAH	2.49	0.95
2:A:401:HXK:CAA	1:C:157:LEU:CD1	2.49	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:283:PHE:CE2	2:D:401:HXK:CAV	2.55	0.88
1:D:283:PHE:CE2	2:D:401:HXK:CBA	2.60	0.85
1:A:165:LYS:NZ	1:C:185:ARG:HH22	1.76	0.83
1:D:283:PHE:HD2	2:D:401:HXK:CAH	1.92	0.79
1:A:190:LYS:HD3	1:A:246:TYR:CD1	2.19	0.78
1:C:265:ASP:CB	4:C:537:HOH:O	2.30	0.78
1:D:190:LYS:HD3	1:D:246:TYR:CD1	2.20	0.77
1:A:262:ILE:CB	4:A:551:HOH:O	2.32	0.77
2:B:401:HXK:CAA	1:D:157:LEU:HD12	2.16	0.76
1:D:283:PHE:HD2	2:D:401:HXK:CAF	2.00	0.75
1:A:221:PRO:HA	1:C:145:LYS:HG3	1.69	0.74
1:A:51:HIS:O	1:A:54:GLU:HG2	1.88	0.73
4:A:545:HOH:O	1:C:242:HIS:HE1	1.69	0.73
1:A:165:LYS:HZ1	1:C:185:ARG:HH22	1.35	0.72
1:C:185:ARG:NH2	1:C:186:ASP:OD1	2.22	0.72
1:C:208:ASP:CG	1:C:209:ALA:H	1.93	0.72
1:C:190:LYS:HD3	1:C:246:TYR:CD1	2.25	0.71
1:C:223:GLN:O	1:C:225:VAL:HG13	1.91	0.71
1:B:165:LYS:NZ	1:D:185:ARG:HH11	1.89	0.70
1:A:100:ILE:N	4:A:502:HOH:O	2.23	0.70
1:C:119:GLN:HE22	2:C:401:HXK:CAS	2.03	0.70
1:B:321:LEU:HD21	1:B:357:HIS:HE1	1.56	0.70
1:A:240:VAL:HG22	1:A:284:GLN:HG2	1.74	0.70
1:D:391:ILE:C	1:D:393:PHE:N	2.41	0.69
1:B:340:ALA:H	1:B:343:LYS:HB2	1.57	0.69
2:D:401:HXK:OAD	4:D:501:HOH:O	2.10	0.69
1:A:319:PHE:HD1	4:A:523:HOH:O	1.74	0.68
1:A:286:HIS:O	1:A:290:LEU:HD12	1.93	0.68
1:A:145:LYS:HG3	1:C:219:ASN:O	1.94	0.68
1:D:283:PHE:HE2	2:D:401:HXK:CAV	2.04	0.67
1:D:321:LEU:HD21	1:D:357:HIS:HE1	1.58	0.67
1:B:190:LYS:HD3	1:B:246:TYR:CD1	2.31	0.66
1:C:290:LEU:O	1:C:292:ILE:N	2.28	0.66
1:A:333:ASN:HD22	1:A:336:LYS:HD3	1.58	0.66
1:A:165:LYS:NZ	1:C:185:ARG:NH2	2.44	0.66
1:A:41:TYR:HB2	1:A:113:TRP:CZ2	2.31	0.65
1:A:145:LYS:HG3	1:C:220:VAL:O	1.97	0.64
1:C:240:VAL:HG22	1:C:284:GLN:HG2	1.78	0.64
1:A:321:LEU:HD21	1:A:357:HIS:HE1	1.63	0.64
1:A:334:TYR:O	4:A:501:HOH:O	2.15	0.63
1:C:313:TRP:HB3	1:C:314:PRO:HD3	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:266:ASN:OD1	1:B:267:LEU:N	2.32	0.63
1:C:338:ASN:HD21	1:C:340:ALA:HB3	1.64	0.63
1:D:306:ILE:HG23	1:D:334:TYR:CD1	2.34	0.62
1:A:321:LEU:HD21	1:A:357:HIS:CE1	2.34	0.62
1:B:286:HIS:O	1:B:290:LEU:HD12	2.00	0.62
1:B:321:LEU:HD21	1:B:357:HIS:CE1	2.34	0.62
2:B:401:HXK:CAS	2:B:401:HXK:CAV	2.78	0.62
1:B:173:ASN:HB2	4:B:506:HOH:O	2.00	0.62
1:D:286:HIS:O	1:D:290:LEU:HD12	1.99	0.62
1:B:165:LYS:HZ1	1:D:185:ARG:HH11	1.47	0.61
1:D:388:LEU:HD13	1:D:392:LEU:HD13	1.82	0.61
1:C:321:LEU:HD21	1:C:357:HIS:HE1	1.66	0.61
1:A:313:TRP:HB3	1:A:314:PRO:HD3	1.82	0.61
1:D:391:ILE:O	1:D:393:PHE:N	2.34	0.61
1:C:342:VAL:C	1:C:344:VAL:H	2.04	0.60
1:A:238:ASN:O	1:A:242:HIS:HD2	1.83	0.60
1:B:240:VAL:HG22	1:B:284:GLN:HG2	1.82	0.60
1:D:283:PHE:HE2	2:D:401:HXK:CAH	2.09	0.60
1:C:283:PHE:CD1	1:C:392:LEU:HG	2.36	0.60
1:D:321:LEU:HD21	1:D:357:HIS:CE1	2.36	0.60
1:C:37:PHE:N	4:C:501:HOH:O	2.34	0.60
2:B:401:HXK:CAA	1:D:157:LEU:CD1	2.79	0.60
1:C:321:LEU:HD21	1:C:357:HIS:CE1	2.36	0.60
1:A:266:ASN:OD1	1:A:267:LEU:N	2.33	0.60
1:C:266:ASN:OD1	1:C:267:LEU:N	2.35	0.60
1:D:313:TRP:HB3	1:D:314:PRO:HD3	1.84	0.60
1:A:51:HIS:CE1	1:A:165:LYS:HE3	2.36	0.59
1:A:217:ASN:ND2	1:C:145:LYS:HD3	2.17	0.59
1:D:342:VAL:O	1:D:344:VAL:N	2.35	0.59
1:B:391:ILE:C	4:B:502:HOH:O	2.39	0.59
1:D:306:ILE:HG23	1:D:334:TYR:CE1	2.38	0.59
1:C:208:ASP:CG	1:C:209:ALA:N	2.56	0.59
1:D:391:ILE:O	1:D:392:LEU:C	2.40	0.59
1:D:266:ASN:OD1	1:D:267:LEU:N	2.35	0.58
1:D:238:ASN:O	1:D:242:HIS:HD2	1.86	0.58
1:C:119:GLN:NE2	2:C:401:HXK:CAS	2.67	0.58
1:C:338:ASN:ND2	1:C:340:ALA:HB3	2.18	0.58
1:A:219:ASN:O	1:A:220:VAL:C	2.41	0.57
1:C:168:GLU:O	1:C:172:ARG:HB3	2.03	0.57
1:A:322:CYS:HA	1:A:352:TYR:CE1	2.40	0.57
1:C:290:LEU:C	1:C:292:ILE:H	2.07	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:322:CYS:HA	1:C:352:TYR:CE1	2.40	0.57
1:D:283:PHE:CD2	2:D:401:HXK:CAF	2.82	0.56
1:B:238:ASN:O	1:B:242:HIS:HD2	1.89	0.56
1:B:313:TRP:HB3	1:B:314:PRO:HD3	1.88	0.56
1:A:168:GLU:O	1:A:172:ARG:HB3	2.06	0.55
1:C:209:ALA:N	1:C:211:ARG:N	2.55	0.55
1:D:240:VAL:O	1:D:244:THR:HG22	2.06	0.55
1:A:322:CYS:HA	1:A:352:TYR:CZ	2.41	0.55
1:D:51:HIS:CE1	1:D:165:LYS:HE3	2.42	0.54
1:C:238:ASN:O	1:C:242:HIS:HD2	1.89	0.54
1:B:292:ILE:C	1:B:355:ARG:NH1	2.61	0.54
1:B:168:GLU:O	1:B:172:ARG:HB3	2.08	0.54
1:B:150:LYS:O	1:D:129:MET:HE2	2.07	0.54
1:D:342:VAL:C	1:D:344:VAL:H	2.10	0.54
1:B:324:GLU:N	1:B:325:PRO:HD2	2.23	0.54
1:C:344:VAL:HG23	4:C:517:HOH:O	2.08	0.54
1:D:326:ASP:OD2	1:D:352:TYR:OH	2.27	0.53
1:C:304:SER:O	1:C:307:GLN:HG3	2.08	0.53
1:B:286:HIS:CE1	1:B:290:LEU:HD11	2.44	0.53
4:A:545:HOH:O	1:C:242:HIS:CE1	2.50	0.53
1:D:114:CYS:HB3	1:D:163:ILE:HG23	1.91	0.53
1:B:185:ARG:HH11	1:D:165:LYS:HZ1	1.57	0.53
1:D:324:GLU:N	1:D:325:PRO:HD2	2.23	0.53
1:B:168:GLU:CD	1:D:185:ARG:HH12	2.12	0.53
1:D:168:GLU:O	1:D:172:ARG:HB3	2.08	0.53
1:A:281:GLU:C	1:A:283:PHE:H	2.12	0.53
1:B:158:LEU:HB2	1:D:192:ILE:HG21	1.92	0.52
1:C:373:ILE:O	1:C:376:LEU:HB2	2.08	0.52
1:C:212:GLU:CG	1:C:213:ILE:N	2.72	0.52
1:B:111:LEU:O	1:B:115:ILE:HG12	2.09	0.52
1:C:222:GLU:O	1:C:224:PRO:N	2.43	0.51
1:A:375:GLU:N	4:A:506:HOH:O	2.42	0.51
1:D:283:PHE:CD2	2:D:401:HXK:CBA	2.89	0.50
1:A:333:ASN:ND2	1:A:336:LYS:HD3	2.26	0.50
1:A:214:ASP:OD1	1:A:216:ASN:N	2.38	0.50
1:B:147:VAL:O	1:B:151:ASN:HB2	2.11	0.50
1:B:290:LEU:O	1:B:292:ILE:N	2.44	0.50
1:C:219:ASN:O	1:C:220:VAL:O	2.30	0.50
1:B:150:LYS:O	1:D:129:MET:CE	2.60	0.50
1:D:286:HIS:CE1	1:D:290:LEU:HD11	2.47	0.50
1:C:283:PHE:CZ	1:C:392:LEU:HG	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:VAL:O	1:B:244:THR:HG22	2.12	0.49
1:D:205:LYS:N	4:D:504:HOH:O	2.38	0.49
1:D:240:VAL:HG22	1:D:284:GLN:HG2	1.95	0.49
1:C:51:HIS:CE1	1:C:165:LYS:HE3	2.48	0.49
1:D:41:TYR:HB2	1:D:113:TRP:CZ2	2.47	0.49
1:A:213:ILE:N	4:A:508:HOH:O	2.46	0.48
1:C:286:HIS:CE1	1:C:290:LEU:HD11	2.48	0.48
4:A:542:HOH:O	1:C:67:LYS:HD3	2.11	0.48
1:A:168:GLU:OE2	1:C:185:ARG:NH1	2.46	0.48
1:B:168:GLU:HG3	4:B:505:HOH:O	2.14	0.48
1:C:286:HIS:O	1:C:290:LEU:HD12	2.13	0.48
1:D:205:LYS:HG3	1:D:206:TYR:HD2	1.79	0.48
1:A:136:ARG:HE	2:A:401:HXK:CAF	2.27	0.48
1:D:233:PHE:HE1	1:D:313:TRP:CD1	2.31	0.48
1:C:147:VAL:O	1:C:151:ASN:HB2	2.14	0.48
1:D:128:ILE:HD11	1:D:152:ALA:HB1	1.95	0.48
1:D:348:LEU:O	1:D:352:TYR:HD2	1.97	0.48
1:B:205:LYS:HG3	1:B:206:TYR:HD2	1.78	0.47
1:B:80:GLY:CA	2:B:401:HXK:CAF	2.92	0.47
1:C:209:ALA:O	1:C:211:ARG:N	2.47	0.47
1:B:286:HIS:NE2	1:B:290:LEU:HD11	2.29	0.47
1:C:388:LEU:HD13	1:C:392:LEU:HD13	1.97	0.47
1:D:283:PHE:CZ	2:D:401:HXK:OAB	2.68	0.47
1:B:185:ARG:HH11	1:D:165:LYS:NZ	2.12	0.47
1:D:373:ILE:O	1:D:376:LEU:HB2	2.15	0.47
1:B:348:LEU:O	1:B:351:GLN:HB3	2.14	0.46
1:C:333:ASN:HB3	1:C:341:CYS:O	2.15	0.46
1:B:122:PHE:CE2	1:B:191:THR:HG21	2.50	0.46
1:A:217:ASN:ND2	1:C:145:LYS:CD	2.78	0.46
1:A:348:LEU:O	1:A:351:GLN:HB2	2.15	0.46
1:A:111:LEU:O	1:A:115:ILE:HG12	2.16	0.46
1:B:292:ILE:C	1:B:355:ARG:HH12	2.19	0.46
1:C:205:LYS:HG3	1:C:206:TYR:HD2	1.81	0.46
1:B:80:GLY:HA3	2:B:401:HXK:CAF	2.46	0.46
1:C:338:ASN:O	1:C:341:CYS:HB2	2.15	0.46
1:C:69:TYR:CE1	1:C:158:LEU:HD22	2.51	0.46
1:C:107:LYS:HE2	1:C:174:GLU:OE2	2.16	0.46
1:C:240:VAL:HG11	1:C:281:GLU:HA	1.98	0.46
1:B:41:TYR:HB2	1:B:113:TRP:CZ2	2.51	0.45
1:B:281:GLU:C	1:B:283:PHE:H	2.19	0.45
1:B:340:ALA:N	1:B:343:LYS:HB2	2.27	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:324:GLU:N	1:C:325:PRO:HD2	2.31	0.45
1:A:322:CYS:HA	1:A:352:TYR:OH	2.16	0.45
1:B:89:ILE:HG23	1:B:105:TRP:HZ3	1.81	0.45
1:A:281:GLU:C	1:A:283:PHE:N	2.70	0.45
1:C:342:VAL:C	1:C:344:VAL:N	2.70	0.45
1:A:107:LYS:HE2	1:A:174:GLU:OE2	2.17	0.45
1:B:290:LEU:C	1:B:292:ILE:N	2.69	0.45
1:B:340:ALA:HA	1:B:343:LYS:N	2.32	0.45
1:A:158:LEU:O	1:A:158:LEU:HG	2.13	0.45
1:B:197:LEU:HB3	1:B:239:ILE:HG12	1.99	0.45
1:D:197:LEU:HB3	1:D:239:ILE:HG12	1.99	0.45
1:A:145:LYS:HE2	1:C:218:ILE:O	2.18	0.44
2:B:401:HXK:OAC	2:B:401:HXK:SAU	2.75	0.44
1:D:158:LEU:O	1:D:158:LEU:HG	2.09	0.44
1:D:388:LEU:C	1:D:388:LEU:HD12	2.37	0.44
1:A:145:LYS:HD3	1:A:145:LYS:HA	1.69	0.44
1:A:154:ASN:HB2	1:C:129:MET:HE1	1.98	0.44
1:D:86:ILE:H	1:D:86:ILE:HG13	1.70	0.44
1:B:239:ILE:HD12	1:B:311:LEU:HD23	1.99	0.44
1:C:281:GLU:C	1:C:283:PHE:H	2.20	0.44
1:C:240:VAL:O	1:C:244:THR:HG22	2.18	0.44
1:D:69:TYR:CE1	1:D:158:LEU:HD22	2.53	0.44
1:D:227:ASP:OD1	1:D:227:ASP:C	2.56	0.44
1:B:233:PHE:HE1	1:B:313:TRP:CD1	2.36	0.43
1:C:223:GLN:O	1:C:224:PRO:C	2.56	0.43
1:C:326:ASP:OD2	1:C:352:TYR:OH	2.36	0.43
1:D:107:LYS:HE2	1:D:174:GLU:OE2	2.18	0.43
1:D:128:ILE:HD11	1:D:152:ALA:CB	2.48	0.43
1:D:111:LEU:O	1:D:115:ILE:HG12	2.17	0.43
1:D:326:ASP:OD1	1:D:329:LYS:HE3	2.19	0.43
1:D:329:LYS:O	1:D:333:ASN:ND2	2.52	0.43
1:C:222:GLU:O	1:C:223:GLN:C	2.56	0.43
1:A:158:LEU:HB2	1:C:192:ILE:HG21	1.99	0.43
1:C:196:HIS:HE1	1:C:206:TYR:OH	2.01	0.43
1:C:391:ILE:O	1:C:392:LEU:C	2.57	0.43
1:A:165:LYS:HZ3	1:C:185:ARG:HH22	1.62	0.43
1:A:223:GLN:HA	1:A:224:PRO:HD3	1.76	0.43
1:C:122:PHE:CE2	1:C:191:THR:HG21	2.54	0.43
1:D:219:ASN:CA	4:D:539:HOH:O	2.67	0.43
1:D:340:ALA:HA	1:D:343:LYS:HB2	2.00	0.43
1:A:205:LYS:HG3	1:A:206:TYR:HD2	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:306:ILE:O	1:D:335:GLY:HA2	2.19	0.43
1:B:338:ASN:N	4:B:504:HOH:O	2.51	0.43
1:D:219:ASN:HB2	1:D:220:VAL:H	1.10	0.43
1:D:391:ILE:C	1:D:393:PHE:H	2.19	0.43
1:A:173:ASN:ND2	4:A:511:HOH:O	2.51	0.42
1:B:165:LYS:HZ3	1:D:185:ARG:HH11	1.67	0.42
1:D:283:PHE:CE2	2:D:401:HXK:OAB	2.72	0.42
1:C:111:LEU:O	1:C:115:ILE:HG12	2.19	0.42
1:C:217:ASN:OD1	1:C:217:ASN:C	2.58	0.42
1:D:281:GLU:C	1:D:283:PHE:H	2.22	0.42
1:D:323:SER:OG	1:D:325:PRO:HG2	2.19	0.42
1:A:281:GLU:O	1:A:283:PHE:N	2.52	0.42
1:A:219:ASN:O	1:A:220:VAL:O	2.37	0.42
1:A:214:ASP:OD1	1:A:214:ASP:C	2.58	0.42
1:D:205:LYS:HG3	1:D:206:TYR:CD2	2.55	0.42
1:B:340:ALA:H	1:B:343:LYS:CB	2.29	0.42
1:C:310:LYS:HA	1:C:310:LYS:HD2	1.92	0.42
1:A:36:PHE:O	1:A:39:ASN:HB2	2.19	0.42
1:A:332:LYS:O	1:A:332:LYS:HG2	2.20	0.42
1:C:84:ARG:HB3	1:C:115:ILE:HG21	2.01	0.42
1:D:147:VAL:O	1:D:151:ASN:HB2	2.20	0.42
1:D:310:LYS:HD2	1:D:310:LYS:HA	1.87	0.42
1:B:388:LEU:HD12	1:B:388:LEU:C	2.40	0.41
1:D:190:LYS:HE2	1:D:246:TYR:HE1	1.85	0.41
1:B:235:VAL:O	1:B:238:ASN:HB2	2.19	0.41
1:D:219:ASN:HA	4:D:539:HOH:O	2.20	0.41
1:A:235:VAL:O	1:A:238:ASN:HB2	2.19	0.41
1:B:227:ASP:C	1:B:227:ASP:OD1	2.59	0.41
1:C:197:LEU:HB3	1:C:239:ILE:HG12	2.01	0.41
1:B:146:ASP:HB2	1:D:213:ILE:HG23	2.02	0.41
1:A:240:VAL:O	1:A:244:THR:HG22	2.21	0.41
1:C:143:LEU:HD12	1:C:143:LEU:HA	1.93	0.41
1:C:205:LYS:HG3	1:C:206:TYR:CD2	2.55	0.41
1:A:129:MET:HE3	1:C:150:LYS:O	2.21	0.41
1:D:190:LYS:HE2	1:D:246:TYR:CE1	2.56	0.41
1:D:216:ASN:OD1	4:D:502:HOH:O	2.22	0.41
1:A:170:TYR:C	1:A:171:LEU:HD12	2.41	0.41
1:B:168:GLU:CD	1:D:185:ARG:NH1	2.75	0.41
1:B:310:LYS:HD2	1:B:310:LYS:HA	1.90	0.41
1:B:349:TYR:C	1:B:351:GLN:H	2.23	0.41
1:C:290:LEU:C	1:C:292:ILE:N	2.73	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:ILE:CB	1:B:355:ARG:HH12	2.34	0.40
1:B:312:THR:HB	1:B:314:PRO:HD2	2.02	0.40
1:D:190:LYS:CE	1:D:246:TYR:CE1	3.04	0.40
1:C:344:VAL:CG2	4:C:517:HOH:O	2.68	0.40
1:D:348:LEU:O	1:D:351:GLN:HB3	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	301/375 (80%)	286 (95%)	15 (5%)	0	100	100
1	B	310/375 (83%)	295 (95%)	13 (4%)	2 (1%)	25	50
1	C	318/375 (85%)	292 (92%)	20 (6%)	6 (2%)	8	20
1	D	320/375 (85%)	300 (94%)	17 (5%)	3 (1%)	17	40
All	All	1249/1500 (83%)	1173 (94%)	65 (5%)	11 (1%)	17	40

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	344	VAL
1	C	291	ASP
1	D	343	LYS
1	D	392	LEU
1	B	291	ASP
1	C	221	PRO
1	D	220	VAL
1	C	343	LYS
1	C	220	VAL
1	C	223	GLN

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Mol	Chain	Res	Type
1	C	224	PRO

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	261/339 (77%)	249 (95%)	12 (5%)	27	54
1	B	258/339 (76%)	244 (95%)	14 (5%)	22	47
1	C	268/339 (79%)	257 (96%)	11 (4%)	30	59
1	D	268/339 (79%)	256 (96%)	12 (4%)	27	55
All	All	1055/1356 (78%)	1006 (95%)	49 (5%)	27	54

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	PHE
1	A	56	SER
1	A	131	LYS
1	A	143	LEU
1	A	172	ARG
1	A	207	SER
1	A	224	PRO
1	A	233	PHE
1	A	328	ILE
1	A	333	ASN
1	A	348	LEU
1	A	388	LEU
1	B	56	SER
1	B	133	GLU
1	B	143	LEU
1	B	158	LEU
1	B	172	ARG
1	B	204	ASP
1	B	226	ILE
1	B	233	PHE

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Mol	Chain	Res	Type
1	B	247	TYR
1	B	328	ILE
1	B	344	VAL
1	B	351	GLN
1	B	355	ARG
1	B	388	LEU
1	C	143	LEU
1	C	172	ARG
1	C	185	ARG
1	C	212	GLU
1	C	219	ASN
1	C	226	ILE
1	C	233	PHE
1	C	328	ILE
1	C	342	VAL
1	C	355	ARG
1	C	388	LEU
1	D	143	LEU
1	D	172	ARG
1	D	204	ASP
1	D	226	ILE
1	D	233	PHE
1	D	247	TYR
1	D	328	ILE
1	D	351	GLN
1	D	355	ARG
1	D	388	LEU
1	D	392	LEU
1	D	393	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	173	ASN
1	A	196	HIS
1	A	238	ASN
1	A	242	HIS
1	A	333	ASN
1	A	357	HIS
1	B	196	HIS
1	B	242	HIS
1	B	286	HIS

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Mol	Chain	Res	Type
1	B	338	ASN
1	B	357	HIS
1	B	374	ASN
1	C	196	HIS
1	C	242	HIS
1	C	286	HIS
1	C	338	ASN
1	C	357	HIS
1	D	196	HIS
1	D	216	ASN
1	D	242	HIS
1	D	337	ASN
1	D	357	HIS

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	SO4	D	402	-	4,4,4	0.22	0	6,6,6	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	HXK	B	401	-	28,28,28	0.89	0	35,35,35	1.03	2 (5%)
2	HXK	C	401	-	28,28,28	1.09	1 (3%)	35,35,35	1.32	4 (11%)
2	HXK	D	401	-	28,28,28	1.00	1 (3%)	35,35,35	0.91	1 (2%)
2	HXK	A	401	-	28,28,28	1.32	2 (7%)	35,35,35	1.05	3 (8%)

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
2	HXK	C	401	-	-	13/18/18/18	0/2/2/2
2	HXK	A	401	-	-	7/18/18/18	0/2/2/2
2	HXK	D	401	-	-	12/18/18/18	0/2/2/2
2	HXK	B	401	-	-	13/18/18/18	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	HXK	CBA-CAV	4.88	1.60	1.49
2	D	401	HXK	CAZ-SAU	-2.75	1.73	1.77
2	C	401	HXK	CAZ-SAU	-2.08	1.74	1.77
2	A	401	HXK	CAZ-SAU	-2.02	1.74	1.77

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	401	HXK	CAS-SAU-CAZ	-4.19	95.14	102.73
2	B	401	HXK	CAZ-CBA-CAV	-3.22	117.49	122.06
2	A	401	HXK	CAS-SAU-CAZ	-2.72	97.79	102.73
2	D	401	HXK	OAC-CAV-CBA	-2.35	116.24	121.94
2	C	401	HXK	CBA-CAZ-SAU	2.23	122.97	120.33
2	C	401	HXK	CAG-CAZ-SAU	-2.22	116.04	121.46
2	A	401	HXK	CAG-CAZ-SAU	-2.21	116.06	121.46
2	B	401	HXK	OAC-CAV-CBA	-2.09	116.85	121.94
2	C	401	HXK	CAX-CAS-SAU	2.07	118.74	110.62
2	A	401	HXK	CBA-CAZ-SAU	2.03	122.74	120.33

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	401	HXK	CAX-CAS-SAU-CAZ
2	B	401	HXK	CBA-CAZ-SAU-CAS
2	C	401	HXK	CAJ-CAY-OAT-CAR
2	C	401	HXK	CAK-CAY-OAT-CAR
2	B	401	HXK	CAK-CAY-OAT-CAR
2	B	401	HXK	CAJ-CAY-OAT-CAR
2	A	401	HXK	CAO-CAP-CAQ-CAR
2	A	401	HXK	CAK-CAY-OAT-CAR
2	B	401	HXK	CAO-CAP-CAQ-CAR
2	A	401	HXK	CAJ-CAY-OAT-CAR
2	D	401	HXK	CAN-CAO-CAP-CAQ
2	B	401	HXK	CAG-CAZ-SAU-CAS
2	C	401	HXK	CAQ-CAR-OAT-CAY
2	C	401	HXK	CAM-CAN-CAO-CAP
2	D	401	HXK	CAK-CAY-OAT-CAR
2	C	401	HXK	SAU-CAS-CAX-CAI
2	D	401	HXK	CAJ-CAY-OAT-CAR
2	B	401	HXK	CAQ-CAR-OAT-CAY
2	A	401	HXK	CAA-CAL-CAM-CAN
2	C	401	HXK	CAA-CAL-CAM-CAN
2	C	401	HXK	CAP-CAQ-CAR-OAT
2	C	401	HXK	SAU-CAS-CAX-CAK
2	D	401	HXK	CAA-CAL-CAM-CAN
2	B	401	HXK	CAA-CAL-CAM-CAN
2	A	401	HXK	CAQ-CAR-OAT-CAY
2	D	401	HXK	CAL-CAM-CAN-CAO
2	B	401	HXK	CAL-CAM-CAN-CAO
2	D	401	HXK	CAM-CAN-CAO-CAP
2	B	401	HXK	SAU-CAS-CAX-CAK
2	D	401	HXK	OAC-CAV-CBA-CAZ
2	B	401	HXK	SAU-CAS-CAX-CAI
2	A	401	HXK	CAX-CAS-SAU-CAZ
2	C	401	HXK	CBA-CAZ-SAU-CAS
2	C	401	HXK	OAC-CAV-CBA-CAH
2	A	401	HXK	CAL-CAM-CAN-CAO
2	C	401	HXK	CAN-CAO-CAP-CAQ
2	C	401	HXK	OAB-CAV-CBA-CAH
2	B	401	HXK	CAM-CAN-CAO-CAP
2	D	401	HXK	OAB-CAV-CBA-CAZ
2	D	401	HXK	OAC-CAV-CBA-CAH
2	D	401	HXK	OAB-CAV-CBA-CAH
2	B	401	HXK	CAX-CAS-SAU-CAZ
2	D	401	HXK	SAU-CAS-CAX-CAK

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Mol	Chain	Res	Type	Atoms
2	D	401	HXK	SAU-CAS-CAX-CAI
2	B	401	HXK	OAB-CAV-CBA-CAH

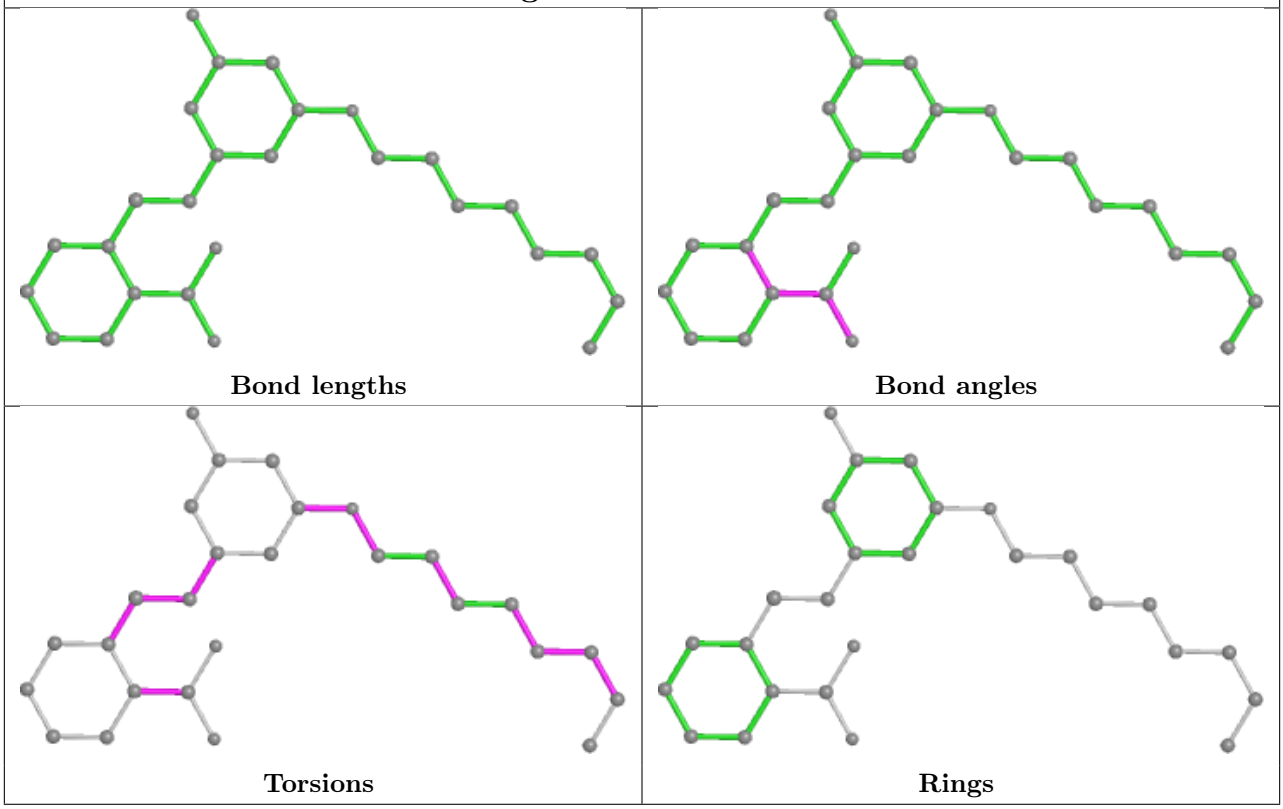
There are no ring outliers.

4 monomers are involved in 24 short contacts:

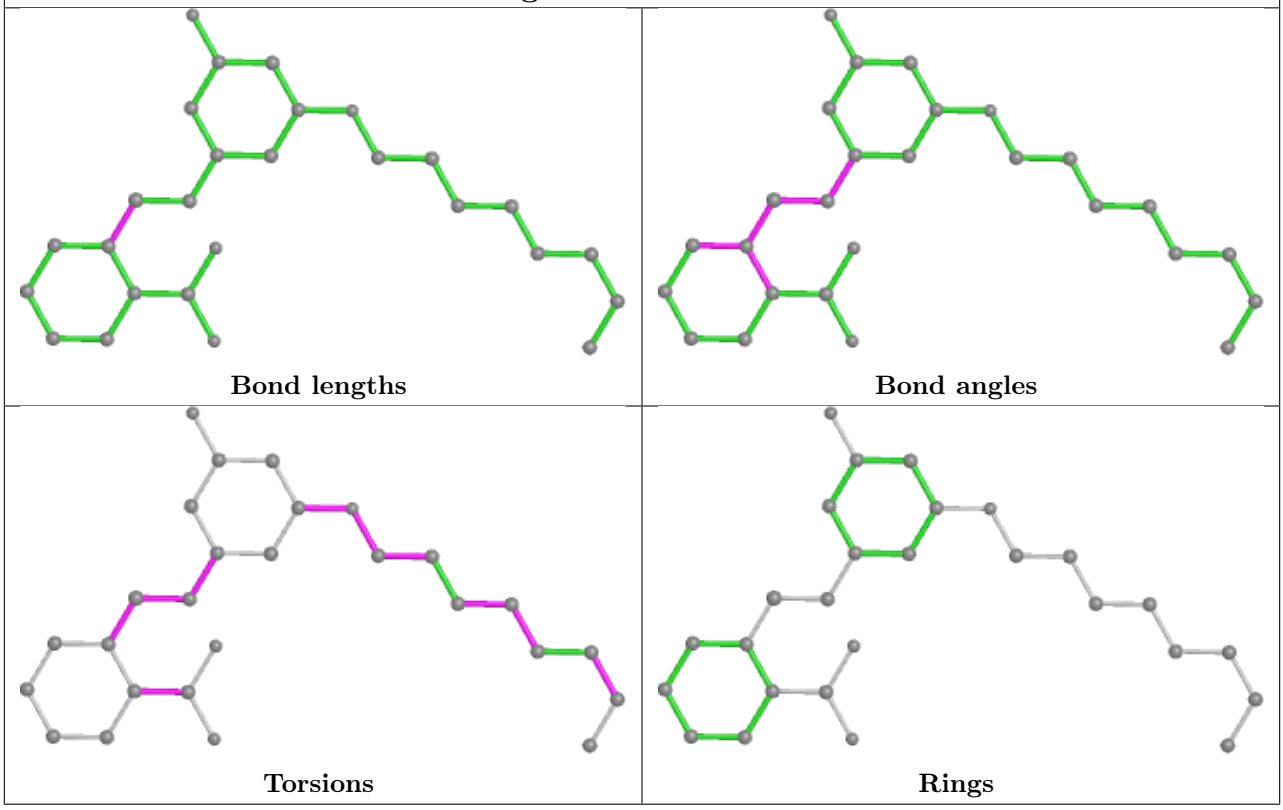
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	HXK	6	0
2	C	401	HXK	2	0
2	D	401	HXK	13	0
2	A	401	HXK	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.

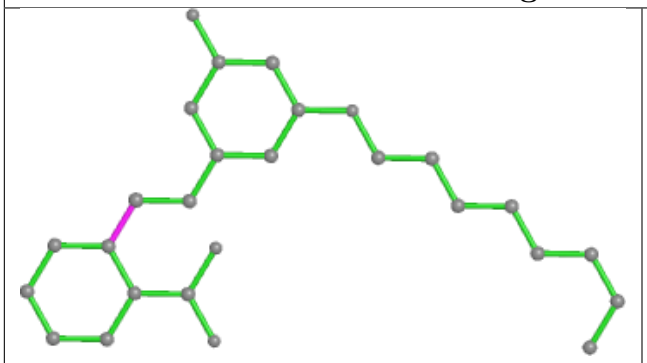
Ligand HXK B 401



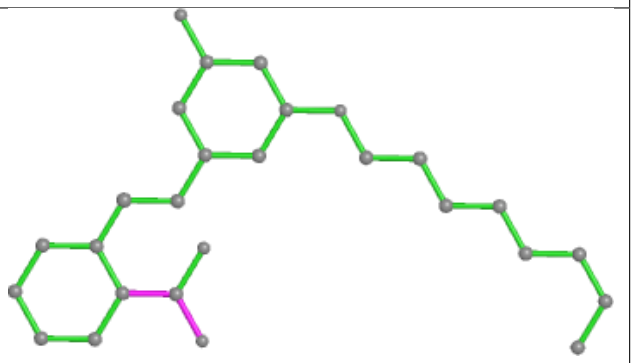
Ligand HXK C 401



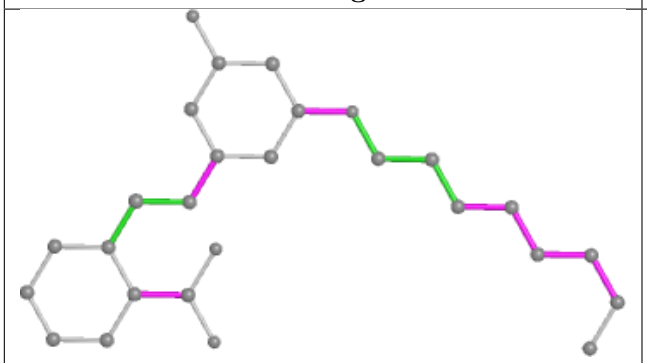
Ligand HXK D 401



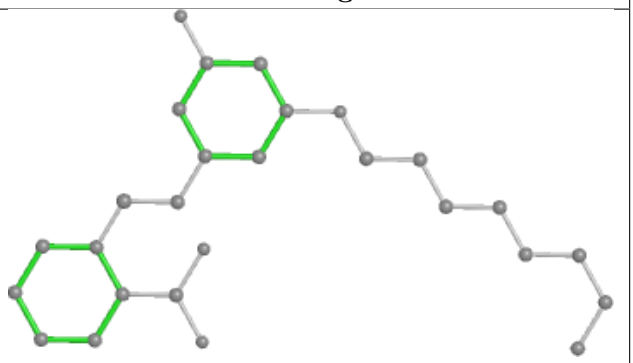
Bond lengths



Bond angles

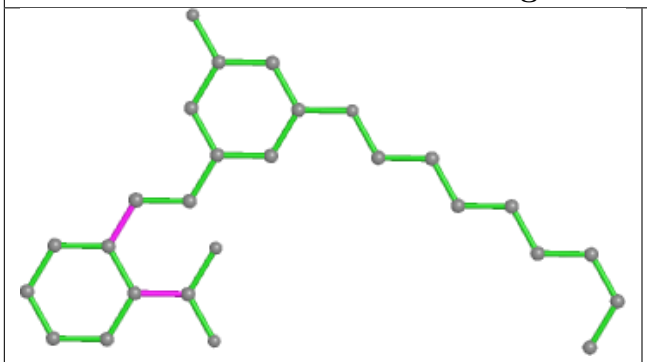


Torsions

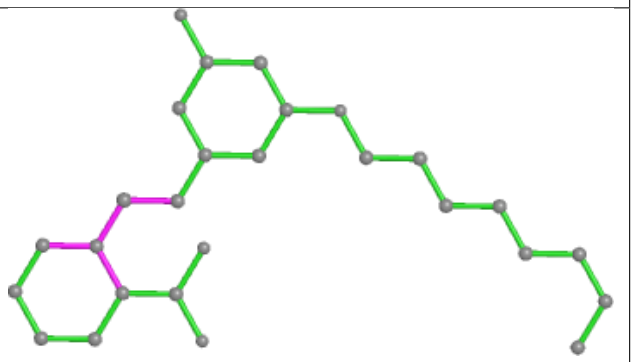


Rings

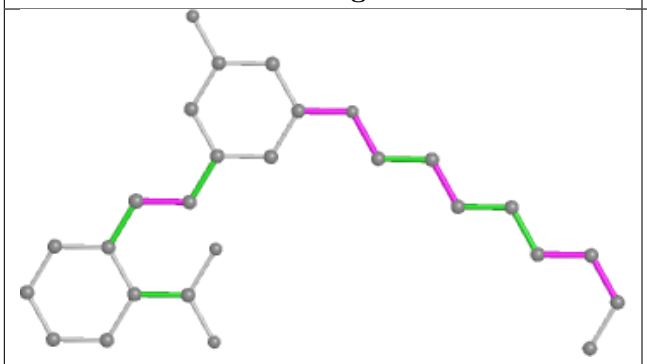
Ligand HXK A 401



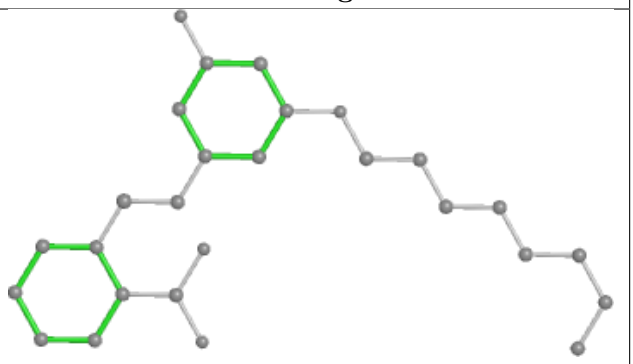
Bond lengths



Bond angles



Torsions



Rings

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	321/375 (85%)	0.38	15 (4%)	31	30	21, 49, 70, 80	0
1	B	326/375 (86%)	0.10	11 (3%)	45	45	21, 49, 72, 99	0
1	C	334/375 (89%)	0.32	17 (5%)	28	26	21, 51, 71, 80	0
1	D	334/375 (89%)	0.12	8 (2%)	59	60	21, 50, 71, 80	0
All	All	1315/1500 (87%)	0.23	51 (3%)	39	38	21, 50, 71, 99	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	283	PHE	4.4
1	A	388	LEU	4.1
1	A	345	ILE	4.1
1	C	224	PRO	3.6
1	A	325	PRO	3.5
1	C	137	ASN	3.5
1	C	100	ILE	3.3
1	B	219	ASN	3.3
1	C	102	SER	3.2
1	A	358	TYR	3.2
1	D	219	ASN	3.2
1	A	387	TYR	3.1
1	B	102	SER	3.1
1	D	137	ASN	3.0
1	A	344	VAL	3.0
1	B	357	HIS	2.8
1	B	224	PRO	2.8
1	C	219	ASN	2.8
1	D	206	TYR	2.8
1	C	349	TYR	2.7
1	D	63	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	283	PHE	2.6
1	A	361	TYR	2.6
1	A	328	ILE	2.5
1	C	61	ILE	2.5
1	C	339	LEU	2.5
1	B	290	LEU	2.5
1	B	100	ILE	2.5
1	A	90	LEU	2.5
1	C	143	LEU	2.4
1	B	226	ILE	2.4
1	A	359	GLU	2.3
1	A	283	PHE	2.3
1	C	358	TYR	2.3
1	B	283	PHE	2.3
1	C	322	CYS	2.3
1	C	220	VAL	2.2
1	D	340	ALA	2.2
1	C	333	ASN	2.2
1	A	386	LYS	2.1
1	D	233	PHE	2.1
1	C	334	TYR	2.1
1	A	333	ASN	2.1
1	C	355	ARG	2.1
1	A	233	PHE	2.1
1	A	348	LEU	2.1
1	B	206	TYR	2.1
1	B	387	TYR	2.1
1	D	290	LEU	2.0
1	C	289	TYR	2.0
1	B	216	ASN	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)

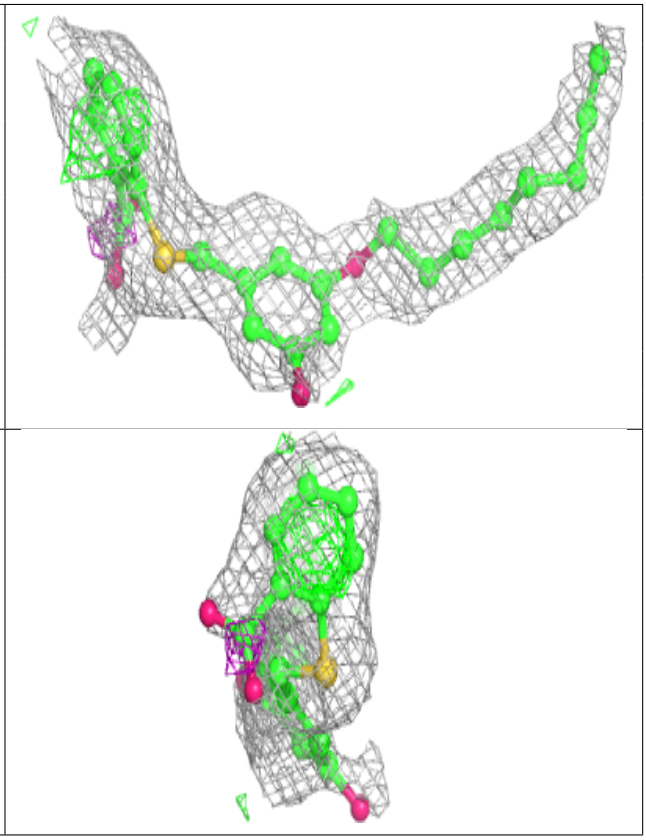
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

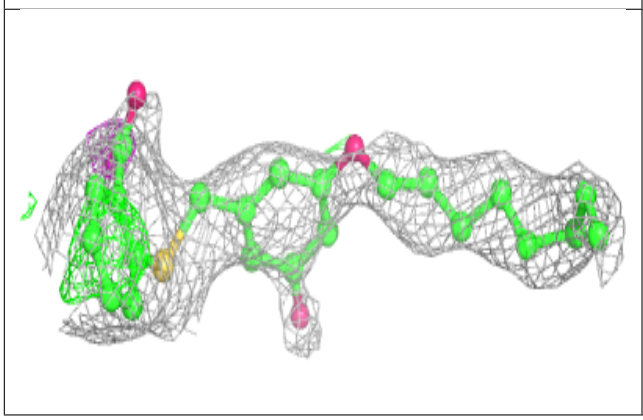
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	HXK	B	401	27/27	0.77	0.29	31,56,62,63	0
2	HXK	C	401	27/27	0.77	0.31	47,63,72,72	0
2	HXK	A	401	27/27	0.81	0.28	34,52,57,59	0
2	HXK	D	401	27/27	0.88	0.24	38,49,53,54	0
3	SO4	D	402	5/5	0.97	0.09	43,43,45,46	0

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 HXK B 401:

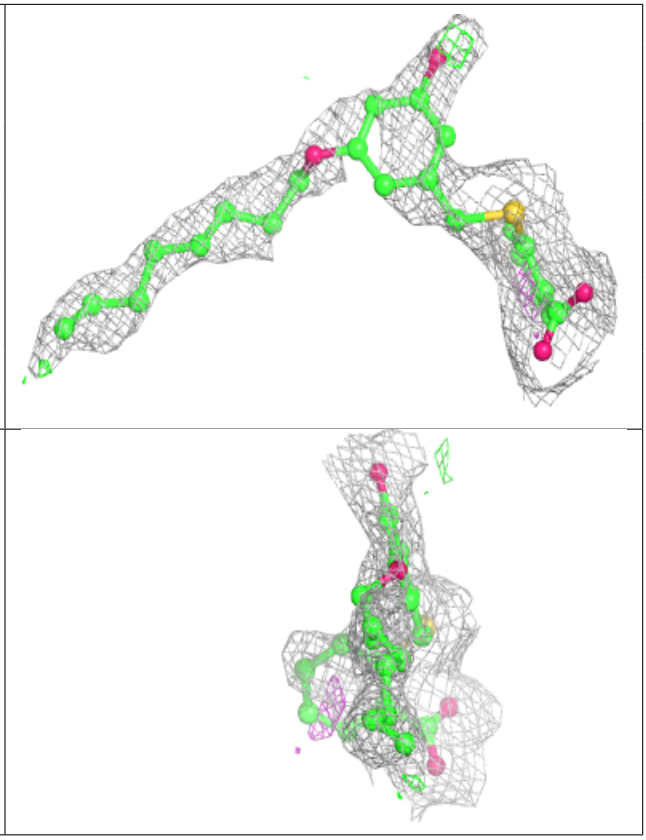
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)





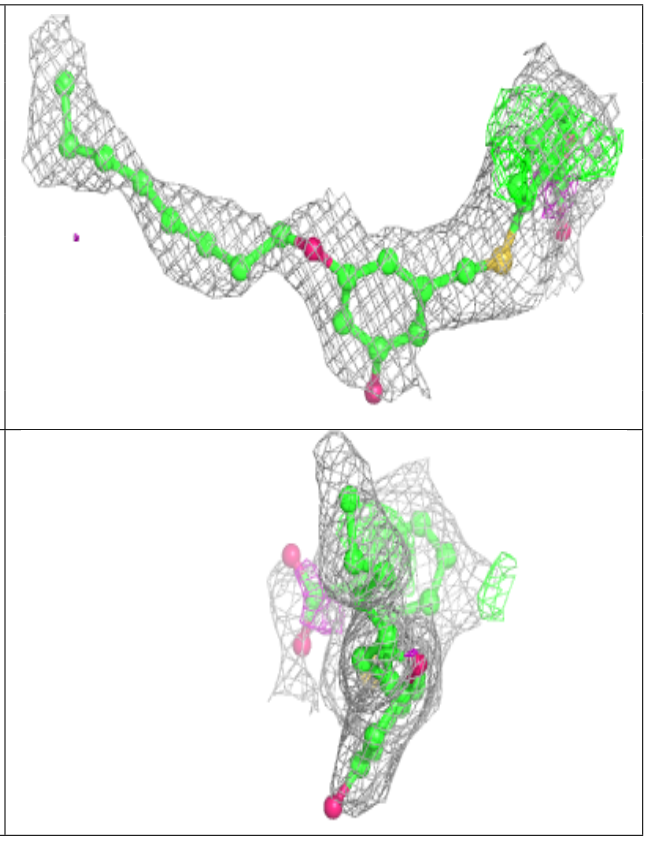
Electron density around HXK C 401:

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



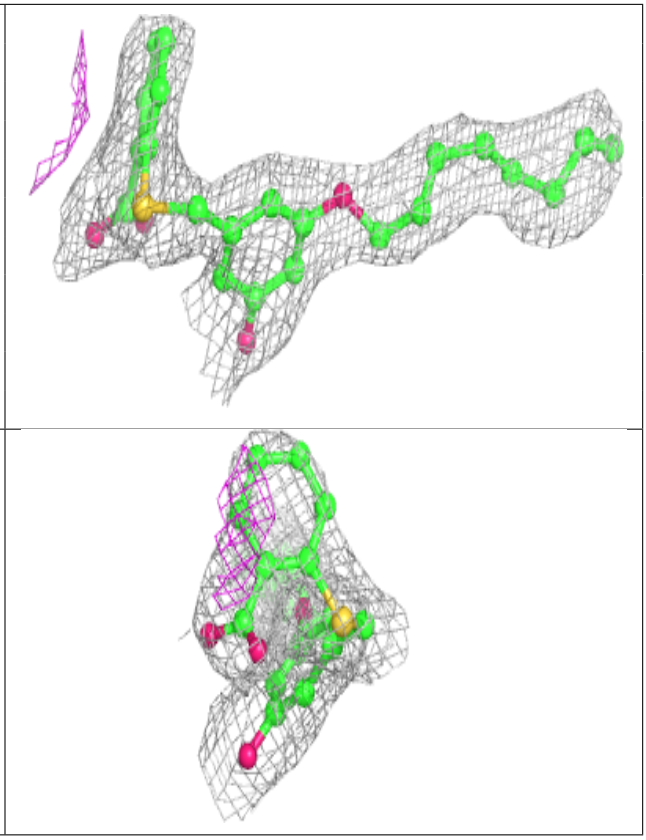
Electron density around HXK A 401:

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



Electron density around HXK D 401:

$2mF_o$-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.