Full wwPDB X-ray Structure Validation Report (i)

Jan 16, 2024 – 03:52 am GMT

PDB ID	:	6YVT
Title	:	HIF prolyl hydroxylase 2 (PHD2/ EGLN1) in complex with MD-253
Authors	:	Chowdhury, R.; Demetriades, M.; Schofield, C.J.
Deposited on	:	2020-04-28
Resolution	:	2.85 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:

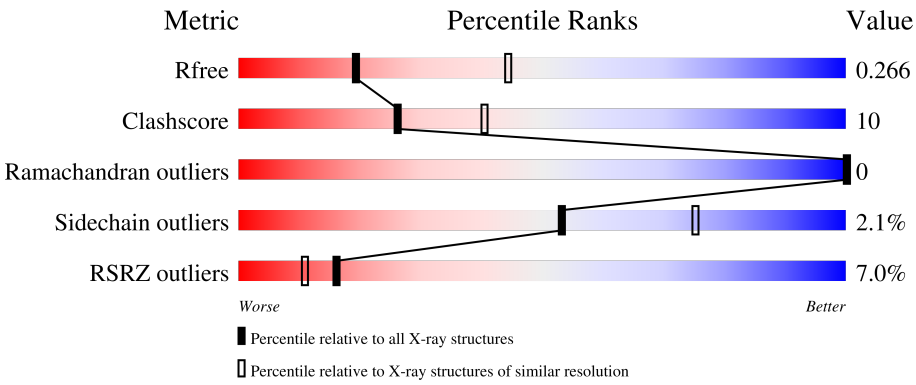
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.85 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	1031 (2.86-2.82)
Clashscore	141614	1078 (2.86-2.82)
Ramachandran outliers	138981	1050 (2.86-2.82)
Sidechain outliers	138945	1051 (2.86-2.82)
RSRZ outliers	127900	1019 (2.86-2.82)

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	252	5% 62% 16% 22%
1	B	252	6% 63% 15% • 21%
1	C	252	5% 62% 15% 22%
1	D	252	6% 60% 18% • 21%
1	E	252	5% 58% 19% 23%

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Mol	Chain	Length	Quality of chain
1	F	252	6% 60% 18% • 22%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 18029 atoms, of which 8671 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 Egl nine homolog 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	197	2877	949	1383	258	276	11	0	0	0
1	B	199	3028	986	1472	272	287	11	0	2	0
1	C	196	2946	960	1432	265	278	11	0	0	0
1	D	199	3007	981	1461	270	284	11	0	0	0
1	E	194	2934	959	1427	260	277	11	0	0	0
1	F	197	2909	954	1398	263	283	11	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	175	GLY	-	expression tag	UNP Q9GZT9
A	176	SER	-	expression tag	UNP Q9GZT9
A	177	HIS	-	expression tag	UNP Q9GZT9
A	178	MET	-	expression tag	UNP Q9GZT9
A	179	ALA	-	expression tag	UNP Q9GZT9
A	180	SER	-	expression tag	UNP Q9GZT9
B	175	GLY	-	expression tag	UNP Q9GZT9
B	176	SER	-	expression tag	UNP Q9GZT9
B	177	HIS	-	expression tag	UNP Q9GZT9
B	178	MET	-	expression tag	UNP Q9GZT9
B	179	ALA	-	expression tag	UNP Q9GZT9
B	180	SER	-	expression tag	UNP Q9GZT9
C	175	GLY	-	expression tag	UNP Q9GZT9
C	176	SER	-	expression tag	UNP Q9GZT9
C	177	HIS	-	expression tag	UNP Q9GZT9
C	178	MET	-	expression tag	UNP Q9GZT9
C	179	ALA	-	expression tag	UNP Q9GZT9

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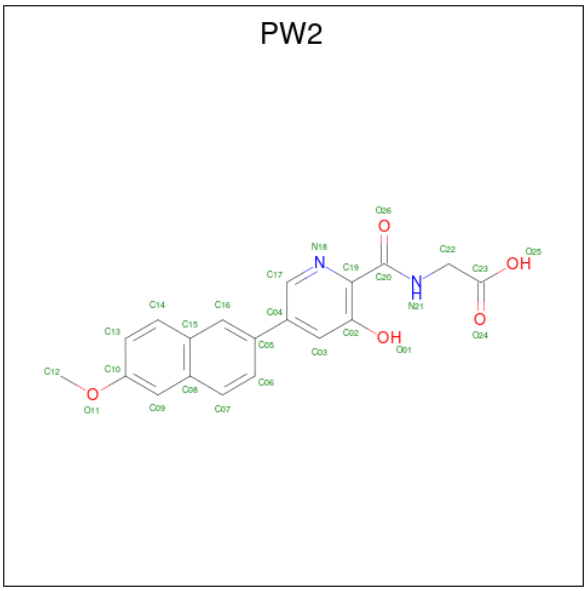
Chain	Residue	Modelled	Actual	Comment	Reference
C	180	SER	-	expression tag	UNP Q9GZT9
D	175	GLY	-	expression tag	UNP Q9GZT9
D	176	SER	-	expression tag	UNP Q9GZT9
D	177	HIS	-	expression tag	UNP Q9GZT9
D	178	MET	-	expression tag	UNP Q9GZT9
D	179	ALA	-	expression tag	UNP Q9GZT9
D	180	SER	-	expression tag	UNP Q9GZT9
E	175	GLY	-	expression tag	UNP Q9GZT9
E	176	SER	-	expression tag	UNP Q9GZT9
E	177	HIS	-	expression tag	UNP Q9GZT9
E	178	MET	-	expression tag	UNP Q9GZT9
E	179	ALA	-	expression tag	UNP Q9GZT9
E	180	SER	-	expression tag	UNP Q9GZT9
F	175	GLY	-	expression tag	UNP Q9GZT9
F	176	SER	-	expression tag	UNP Q9GZT9
F	177	HIS	-	expression tag	UNP Q9GZT9
F	178	MET	-	expression tag	UNP Q9GZT9
F	179	ALA	-	expression tag	UNP Q9GZT9
F	180	SER	-	expression tag	UNP Q9GZT9

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

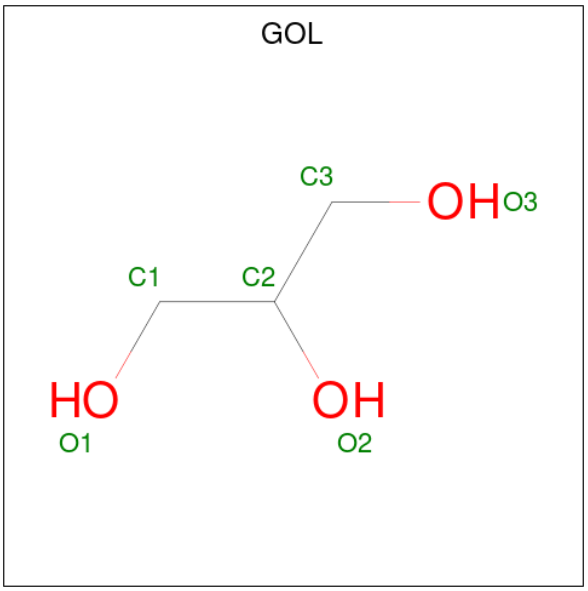
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mn		
2	A	1	1	1	0	0
2	B	1	1	1	0	0
2	C	1	1	1	0	0
2	D	1	1	1	0	0
2	E	1	1	1	0	0
2	F	1	1	1	0	0

- Molecule 3 is 2-[[5-(6-methoxynaphthalen-2-yl)-3-oxidanyl-pyridin-2-yl]carbonylamino]ethanoic acid (three-letter code: PW2) (formula: $C_{19}H_{16}N_2O_5$) (labeled as "Ligand of Interest" by depositor).



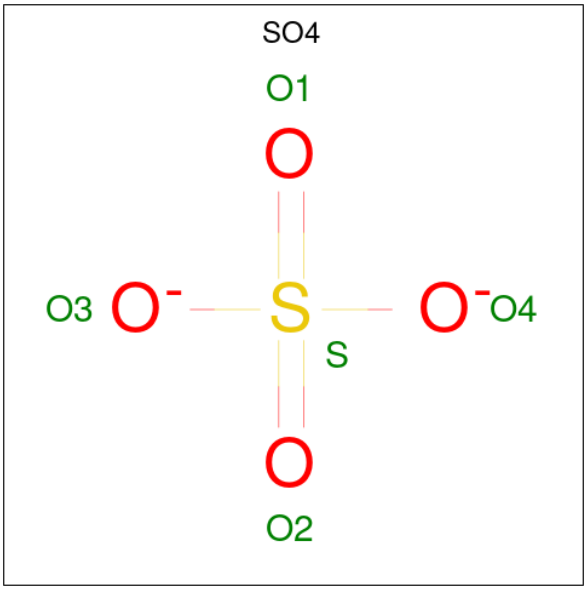
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			41	19	15	2	5		
3	B	1	Total	C	H	N	O	0	0
			41	19	15	2	5		
3	C	1	Total	C	H	N	O	0	0
			41	19	15	2	5		
3	D	1	Total	C	H	N	O	0	0
			41	19	15	2	5		
3	E	1	Total	C	H	N	O	0	0
			41	19	15	2	5		
3	F	1	Total	C	H	N	O	0	0
			41	19	15	2	5		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
4	E	1	14	3	8	3	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 6 is water.

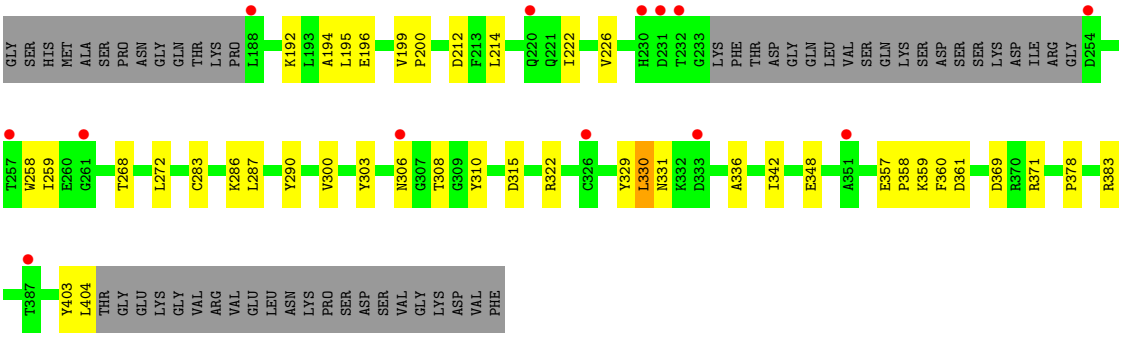
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	6	6	6	0	0
			Total	O		
6	B	7	7	7	0	0
			Total	O		
6	C	9	9	9	0	0
			Total	O		
6	D	13	13	13	0	0
			Total	O		
6	E	14	14	14	0	0
			Total	O		
6	F	8	8	8	0	0

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.

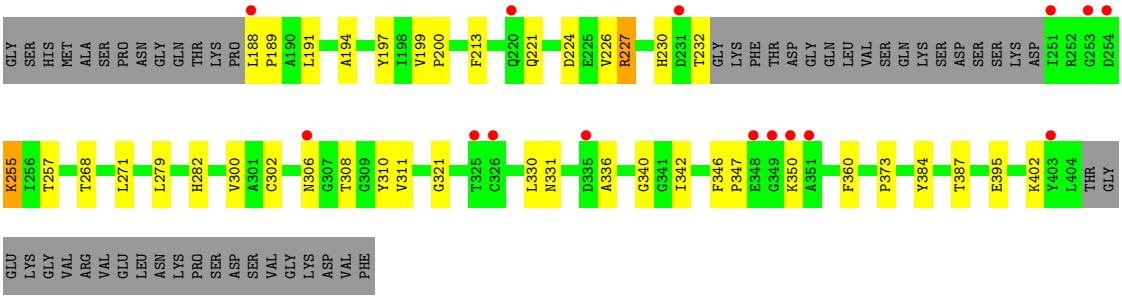
• Molecule 1: Egl nine homolog 1

Chain A:



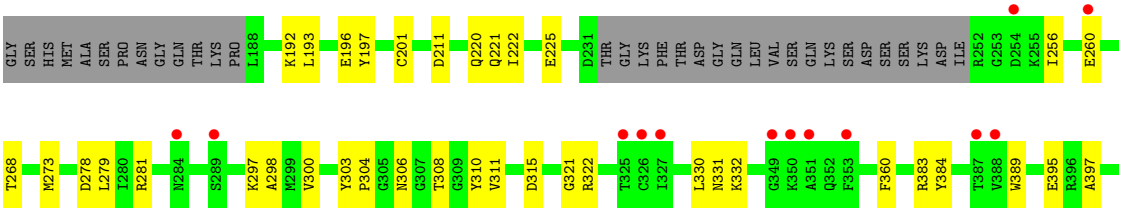
• Molecule 1: Egl nine homolog 1

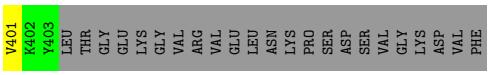
Chain B:



• Molecule 1: Egl nine homolog 1

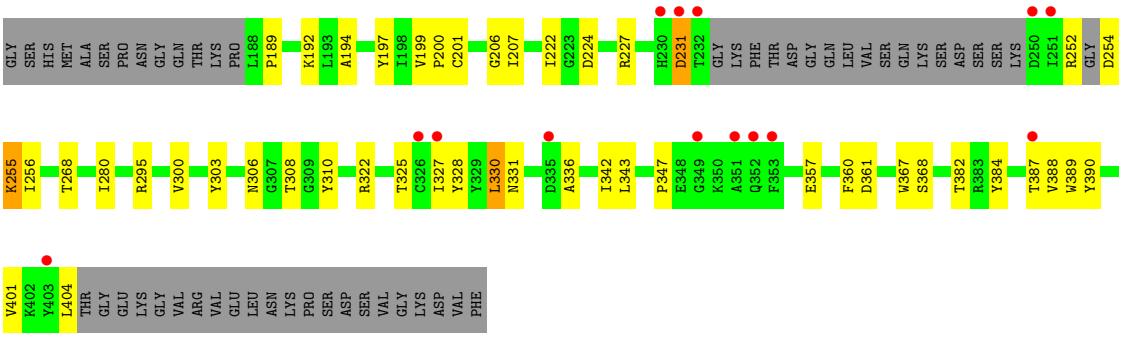
Chain C:





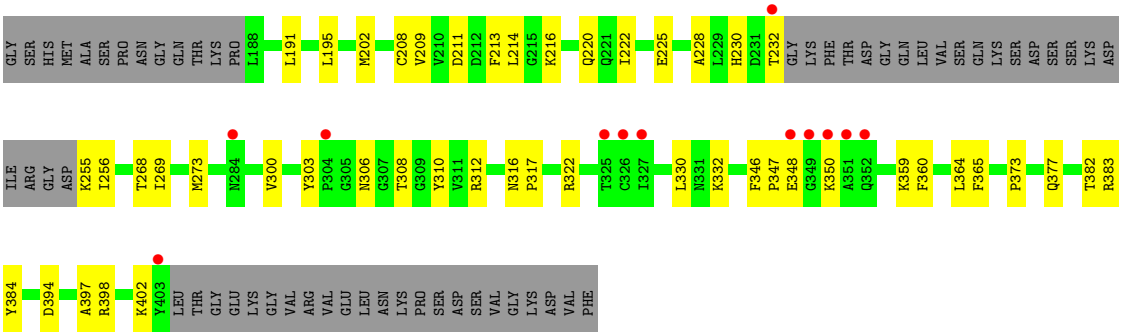
• Molecule 1: Egl nine homolog 1

Chain D:



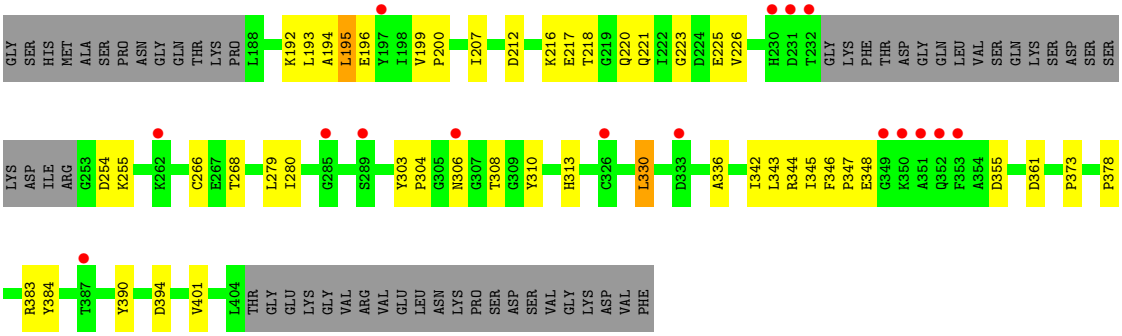
• Molecule 1: Egl nine homolog 1

Chain E:



• Molecule 1: Egl nine homolog 1

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	84.46Å 103.02Å 196.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	23.75 – 2.85 23.75 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.4 (23.75-2.85) 99.5 (23.75-2.85)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.81 (at 2.84Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.237 , 0.266 0.237 , 0.266	Depositor DCC
R_{free} test set	2041 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	48.4	Xtriage
Anisotropy	0.213	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.45 , 57.9	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	0.91	EDS
Total number of atoms	18029	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.45 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.6664e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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, GOL, PW2, MN

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	0.30	0/1530	0.51	0/2082
1	B	0.30	0/1598	0.54	0/2166
1	C	0.30	0/1550	0.52	0/2103
1	D	0.30	0/1582	0.54	0/2145
1	E	0.30	0/1544	0.53	0/2096
1	F	0.30	0/1547	0.52	0/2104
All	All	0.30	0/9351	0.53	0/12696

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1494	1383	1383	30	0
1	B	1556	1472	1467	27	0
1	C	1514	1432	1433	22	0
1	D	1546	1461	1460	39	0
1	E	1507	1427	1432	37	0
1	F	1511	1398	1399	34	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	26	15	0	1	0
3	B	26	15	0	1	0
3	C	26	15	0	3	0
3	D	26	15	0	3	0
3	E	26	15	0	1	0
3	F	26	15	0	1	0
4	E	6	8	8	2	0
5	E	5	0	0	1	0
6	A	6	0	0	0	0
6	B	7	0	0	1	0
6	C	9	0	0	0	0
6	D	13	0	0	1	0
6	E	14	0	0	2	0
6	F	8	0	0	0	0
All	All	9358	8671	8582	186	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 10.

All (186) 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:F:306:ASN:OD1	1:F:308:THR:HG23	1.68	0.94
1:A:336:ALA:HB1	1:A:342:ILE:HD11	1.56	0.85
1:C:306:ASN:OD1	1:C:308:THR:HG23	1.77	0.84
1:D:330:LEU:HD11	1:D:384:TYR:HB2	1.65	0.79
1:D:330:LEU:HD11	1:D:384:TYR:CB	2.14	0.77
1:E:306:ASN:OD1	1:E:308:THR:HG23	1.85	0.77
3:C:502:PW2:C06	1:D:401:VAL:HG22	2.16	0.75
1:C:278:ASP:OD1	1:C:281:ARG:NH1	2.21	0.74
1:F:254:ASP:OD1	1:F:255:LYS:N	2.21	0.73
1:C:256:ILE:HA	1:C:300:VAL:O	1.88	0.73
1:E:211:ASP:N	4:E:503:GOL:O1	2.20	0.73
1:C:220:GLN:OE1	1:C:332:LYS:NZ	2.21	0.73
1:D:207:ILE:HD11	1:D:347:PRO:HG3	1.71	0.72
1:E:230:HIS:ND1	1:E:384:TYR:OH	2.22	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:194:ALA:O	1:F:199:VAL:HG23	1.92	0.69
1:B:321:GLY:N	1:B:395:GLU:OE2	2.27	0.67
1:D:325:THR:HG1	1:D:368:SER:HG	1.37	0.67
1:D:255:LYS:NZ	6:D:601:HOH:O	2.27	0.66
1:E:330:LEU:HD11	1:E:384:TYR:HB2	1.78	0.64
1:D:306:ASN:OD1	1:D:308:THR:HG23	1.98	0.64
1:F:336:ALA:HB1	1:F:342:ILE:HD11	1.80	0.63
1:B:336:ALA:HB1	1:B:342:ILE:HD11	1.79	0.62
1:B:224:ASP:OD1	1:B:227[B]:ARG:NH2	2.33	0.62
1:E:228:ALA:O	1:E:232:THR:HG23	2.00	0.62
1:D:194:ALA:O	1:D:199:VAL:HG23	2.01	0.60
1:D:330:LEU:CD1	1:D:384:TYR:HB2	2.30	0.60
1:F:342:ILE:HG21	1:F:355:ASP:HB3	1.83	0.60
1:D:222:ILE:CG1	1:D:268:THR:HG22	2.31	0.60
1:B:330:LEU:HD11	1:B:384:TYR:CB	2.31	0.60
1:E:222:ILE:HG12	1:E:268:THR:HG22	1.82	0.60
1:F:195:LEU:HD13	1:F:199:VAL:HG21	1.82	0.60
1:C:193:LEU:HD23	1:C:279:LEU:HD21	1.83	0.59
1:D:330:LEU:HD11	1:D:384:TYR:HB3	1.85	0.59
1:E:222:ILE:CG1	1:E:268:THR:HG22	2.33	0.58
1:A:222:ILE:CG1	1:A:268:THR:HG22	2.32	0.58
1:E:330:LEU:HD11	1:E:384:TYR:CB	2.33	0.58
1:B:306:ASN:OD1	1:B:308:THR:HG23	2.03	0.58
1:C:304:PRO:HD2	1:C:308:THR:HG21	1.85	0.57
1:D:343:LEU:HD13	3:D:502:PW2:C22	2.33	0.57
1:E:256:ILE:HA	1:E:300:VAL:O	2.05	0.57
1:F:216:LYS:O	1:F:220:GLN:HG3	2.05	0.57
1:C:321:GLY:N	1:C:395:GLU:OE2	2.33	0.57
1:D:222:ILE:HG12	1:D:268:THR:HG22	1.87	0.56
1:B:330:LEU:HD11	1:B:384:TYR:HB3	1.87	0.56
1:A:329:TYR:C	1:A:330:LEU:HD23	2.26	0.55
1:B:347:PRO:HB2	1:B:350:LYS:HG2	1.89	0.55
1:D:207:ILE:CD1	1:D:347:PRO:HG3	2.36	0.55
1:A:222:ILE:HG13	1:A:268:THR:HG22	1.89	0.54
1:E:303:TYR:HB2	1:E:383:ARG:HB3	1.89	0.54
1:F:303:TYR:HB2	1:F:383:ARG:HB3	1.90	0.54
1:D:224:ASP:OD1	1:D:227:ARG:NH2	2.40	0.54
1:D:336:ALA:HB1	1:D:342:ILE:HG13	1.89	0.54
1:D:231:ASP:C	1:D:231:ASP:OD1	2.47	0.53
1:A:192:LYS:HE2	1:A:196:GLU:OE1	2.09	0.53
1:D:327:ILE:HG12	1:D:387:THR:HG23	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:225:GLU:CD	1:F:268:THR:OG1	2.47	0.53
1:E:359:LYS:NZ	5:E:504:SO4:O1	2.24	0.53
1:B:346:PHE:O	1:B:373:PRO:HD2	2.08	0.53
3:A:502:PW2:N21	3:A:502:PW2:O01	2.42	0.53
1:A:330:LEU:HD23	1:A:330:LEU:N	2.24	0.53
3:B:502:PW2:C06	1:F:401:VAL:HG22	2.39	0.53
1:A:310:TYR:OH	1:C:397:ALA:O	2.20	0.52
1:B:188:LEU:HB3	1:B:189:PRO:CD	2.39	0.52
1:F:348:GLU:O	1:F:348:GLU:CG	2.58	0.52
1:E:211:ASP:HB2	4:E:503:GOL:O1	2.09	0.52
1:E:330:LEU:CD1	1:E:384:TYR:HB2	2.41	0.51
1:B:194:ALA:O	1:B:199:VAL:HG23	2.11	0.51
1:A:283:CYS:HB3	1:A:287:LEU:HG	1.93	0.51
1:C:315:ASP:HA	1:D:401:VAL:HG11	1.92	0.51
1:B:311:VAL:HG11	1:F:394:ASP:OD1	2.11	0.51
1:A:357:GLU:HB2	1:A:359:LYS:HG3	1.94	0.50
1:E:397:ALA:O	1:F:310:TYR:OH	2.27	0.50
1:E:191:LEU:HD11	1:E:195:LEU:CD1	2.42	0.50
1:B:331:ASN:O	1:B:360:PHE:HB2	2.12	0.49
1:E:346:PHE:O	1:E:373:PRO:HD2	2.11	0.49
1:E:202:MET:HE1	1:E:365:PHE:HB2	1.93	0.49
1:D:252:ARG:HD3	1:D:254:ASP:O	2.13	0.49
1:D:331:ASN:O	1:D:360:PHE:HB2	2.12	0.49
1:B:191:LEU:HD12	1:B:282:HIS:HB3	1.94	0.49
1:A:195:LEU:HD21	1:A:286:LYS:HE3	1.94	0.48
1:E:220:GLN:NE2	1:E:332:LYS:NZ	2.61	0.48
1:C:260:GLU:HB2	1:C:297:LYS:HD3	1.95	0.48
1:D:222:ILE:HG13	1:D:268:THR:HG22	1.95	0.48
1:D:252:ARG:CD	1:D:254:ASP:O	2.62	0.48
1:C:222:ILE:HG13	1:C:268:THR:HG22	1.95	0.48
1:D:252:ARG:NH2	1:D:303:TYR:CZ	2.82	0.47
1:E:394:ASP:O	1:E:398:ARG:HG3	2.14	0.47
1:B:336:ALA:O	1:B:340:GLY:N	2.46	0.47
1:A:222:ILE:HG12	1:A:268:THR:HG22	1.96	0.47
1:F:218:THR:HA	1:F:221:GLN:HG3	1.97	0.47
1:F:308:THR:O	1:F:378:PRO:HA	2.14	0.47
1:D:280:ILE:HG13	1:D:390:TYR:CZ	2.50	0.47
1:A:308:THR:O	1:A:378:PRO:HA	2.15	0.46
1:A:226:VAL:HG21	1:A:330:LEU:CD1	2.45	0.46
1:A:310:TYR:OH	1:C:401:VAL:HG23	2.16	0.46
1:E:348:GLU:O	6:E:601:HOH:O	2.21	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:269:ILE:O	1:E:273:MET:HG3	2.15	0.46
1:F:192:LYS:O	1:F:196:GLU:HB2	2.15	0.46
1:F:345:ILE:O	1:F:347:PRO:HD3	2.15	0.46
1:C:192:LYS:HE3	1:C:196:GLU:OE1	2.15	0.46
1:D:189:PRO:HD2	1:D:192:LYS:HB3	1.98	0.46
1:E:225:GLU:CD	1:E:268:THR:HB	2.35	0.46
1:B:336:ALA:CB	1:B:342:ILE:HD11	2.45	0.46
1:B:310:TYR:OH	1:F:401:VAL:HG23	2.14	0.46
1:A:194:ALA:O	1:A:199:VAL:HG23	2.16	0.46
1:A:303:TYR:HB2	1:A:383:ARG:HB3	1.98	0.46
1:F:348:GLU:O	1:F:348:GLU:HG2	2.16	0.46
1:F:343:LEU:HD21	1:F:345:ILE:HD11	1.97	0.45
1:D:303:TYR:OH	3:D:502:PW2:O01	2.27	0.45
1:E:213:PHE:CD2	1:E:214:LEU:HG	2.51	0.45
1:E:216:LYS:O	1:E:220:GLN:HB2	2.16	0.45
1:F:217:GLU:O	1:F:221:GLN:HG3	2.16	0.45
1:E:191:LEU:HD11	1:E:195:LEU:HD12	1.97	0.45
1:F:199:VAL:N	1:F:200:PRO:HD2	2.30	0.45
1:F:225:GLU:OE1	1:F:266:CYS:HA	2.15	0.45
1:C:303:TYR:HB2	1:C:383:ARG:HB3	1.97	0.45
1:E:230:HIS:HA	1:E:255:LYS:CB	2.46	0.45
1:D:303:TYR:O	1:D:382:THR:HA	2.15	0.45
1:D:322:ARG:HD2	1:D:389:TRP:CG	2.52	0.45
1:E:316:ASN:ND2	1:E:322:ARG:O	2.40	0.45
1:F:223:GLY:HA2	1:F:330:LEU:CD1	2.47	0.45
1:A:315:ASP:O	1:A:322:ARG:NH2	2.50	0.45
1:B:387:THR:HG21	6:B:604:HOH:O	2.17	0.45
1:A:369:ASP:OD1	1:A:371:ARG:HB2	2.17	0.45
1:E:202:MET:HE1	1:E:208:CYS:HB3	2.00	0.44
1:C:197:TYR:O	1:C:201:CYS:HB2	2.17	0.44
1:F:344:ARG:NE	1:F:355:ASP:OD1	2.47	0.44
1:D:295:ARG:HD2	1:D:388:VAL:HG11	1.99	0.44
1:A:331:ASN:HD22	1:A:358:PRO:HB2	1.82	0.44
1:B:221:GLN:HB3	1:B:268:THR:HG21	2.00	0.44
1:C:330:LEU:HD11	1:C:384:TYR:HB3	1.99	0.44
1:F:313:HIS:O	1:F:373:PRO:HA	2.18	0.44
1:D:256:ILE:HA	1:D:300:VAL:O	2.18	0.44
1:E:220:GLN:NE2	1:E:332:LYS:HZ1	2.16	0.44
1:B:188:LEU:HB3	1:B:189:PRO:HD2	2.00	0.43
3:D:502:PW2:O01	3:D:502:PW2:N21	2.50	0.43
1:A:199:VAL:HB	1:A:200:PRO:CD	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:273:MET:HE3	1:C:298:ALA:HB2	2.00	0.43
3:C:502:PW2:C07	1:D:404:LEU:HD13	2.48	0.43
1:F:207:ILE:HD11	1:F:345:ILE:HG21	2.00	0.43
1:C:221:GLN:O	1:C:225:GLU:HG3	2.18	0.43
1:C:330:LEU:HD11	1:C:384:TYR:CB	2.48	0.43
1:A:403:TYR:O	1:A:404:LEU:C	2.57	0.43
1:B:336:ALA:HB1	1:B:342:ILE:CD1	2.48	0.43
1:E:377:GLN:OE1	6:E:602:HOH:O	2.21	0.43
1:F:280:ILE:HG13	1:F:390:TYR:CE2	2.54	0.43
1:F:225:GLU:OE1	1:F:268:THR:OG1	2.36	0.43
1:B:402:LYS:HG3	1:E:317:PRO:CG	2.48	0.42
1:B:199:VAL:N	1:B:200:PRO:HD2	2.34	0.42
1:D:206:GLY:HA2	1:D:367:TRP:NE1	2.34	0.42
1:E:303:TYR:O	1:E:382:THR:HA	2.19	0.42
3:F:502:PW2:O01	3:F:502:PW2:N21	2.52	0.42
1:F:346:PHE:O	1:F:373:PRO:HD2	2.20	0.42
1:D:336:ALA:HB2	1:D:357:GLU:HG2	2.01	0.42
1:A:306:ASN:OD1	1:A:308:THR:HG23	2.19	0.42
1:A:331:ASN:O	1:A:360:PHE:HB2	2.20	0.42
1:D:336:ALA:CB	1:D:342:ILE:HG13	2.50	0.42
1:A:214:LEU:HD11	1:A:272:LEU:HD12	2.01	0.42
1:B:213:PHE:CE1	1:B:279:LEU:HD11	2.55	0.42
1:D:197:TYR:O	1:D:201:CYS:HB2	2.20	0.42
1:D:222:ILE:HG13	1:D:268:THR:CG2	2.50	0.42
1:E:347:PRO:HG2	1:E:350:LYS:HE3	2.02	0.41
1:C:331:ASN:O	1:C:360:PHE:HB2	2.20	0.41
3:C:502:PW2:O01	3:C:502:PW2:N21	2.53	0.41
1:A:348:GLU:N	1:A:371:ARG:O	2.50	0.41
1:E:222:ILE:HG13	1:E:268:THR:HG22	2.03	0.41
1:E:209:VAL:HG22	1:E:364:LEU:HD13	2.02	0.41
1:F:193:LEU:HD23	1:F:279:LEU:HD21	2.02	0.41
1:A:258:TRP:O	1:A:259:ILE:HD13	2.20	0.41
1:B:226:VAL:HG13	1:B:300:VAL:CG1	2.51	0.41
1:C:322:ARG:HD2	1:C:389:TRP:CG	2.55	0.41
1:D:328:TYR:OH	1:D:361:ASP:OD1	2.20	0.41
1:F:304:PRO:HD2	1:F:308:THR:HG21	2.03	0.41
1:A:226:VAL:HG13	1:A:300:VAL:HG11	2.02	0.41
1:B:255:LYS:HE3	1:B:302:CYS:HB3	2.03	0.41
1:C:256:ILE:HG13	1:C:300:VAL:O	2.21	0.41
1:A:199:VAL:HB	1:A:200:PRO:HD3	2.03	0.41
1:A:212:ASP:HA	1:A:361:ASP:OD2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:199:VAL:N	1:D:200:PRO:HD2	2.36	0.41
1:F:212:ASP:HA	1:F:361:ASP:OD2	2.21	0.41
1:A:290:TYR:OH	1:A:371:ARG:NH2	2.54	0.40
1:B:226:VAL:HG13	1:B:300:VAL:HG11	2.04	0.40
1:F:226:VAL:CG1	1:F:384:TYR:CD1	3.05	0.40
1:E:216:LYS:HA	1:E:360:PHE:HE2	1.86	0.40
1:E:383:ARG:NH2	3:E:502:PW2:O24	2.44	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	193/252 (77%)	184 (95%)	9 (5%)	0	100	100
1	B	197/252 (78%)	193 (98%)	4 (2%)	0	100	100
1	C	192/252 (76%)	189 (98%)	3 (2%)	0	100	100
1	D	193/252 (77%)	189 (98%)	4 (2%)	0	100	100
1	E	190/252 (75%)	187 (98%)	3 (2%)	0	100	100
1	F	193/252 (77%)	181 (94%)	12 (6%)	0	100	100
All	All	1158/1512 (77%)	1123 (97%)	35 (3%)	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	145/211 (69%)	144 (99%)	1 (1%)	84	91
1	B	156/211 (74%)	149 (96%)	7 (4%)	27	51
1	C	152/211 (72%)	149 (98%)	3 (2%)	55	76
1	D	155/211 (74%)	151 (97%)	4 (3%)	46	70
1	E	153/211 (72%)	150 (98%)	3 (2%)	55	76
1	F	150/211 (71%)	148 (99%)	2 (1%)	69	84
All	All	911/1266 (72%)	891 (98%)	20 (2%)	53	75

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	330	LEU
1	B	197	TYR
1	B	227[A]	ARG
1	B	227[B]	ARG
1	B	230	HIS
1	B	255	LYS
1	B	257	THR
1	B	271	LEU
1	C	211	ASP
1	C	310	TYR
1	C	311	VAL
1	D	231	ASP
1	D	255	LYS
1	D	310	TYR
1	D	330	LEU
1	E	310	TYR
1	E	312	ARG
1	E	402	LYS
1	F	195	LEU
1	F	330	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	220	GLN
1	E	377	GLN

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 14 ligands modelled in this entry, 6 are monoatomic - leaving 8 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	$\#\|Z\| > 2$
5	SO4	E	504	-	4,4,4	0.14	0	6,6,6	0.07	0
3	PW2	D	502	2	28,28,28	1.41	2 (7%)	37,39,39	1.85	8 (21%)
3	PW2	E	502	2	28,28,28	1.51	3 (10%)	37,39,39	1.42	5 (13%)
3	PW2	B	502	2	28,28,28	1.47	3 (10%)	37,39,39	1.61	6 (16%)
4	GOL	E	503	-	5,5,5	0.09	0	5,5,5	0.24	0
3	PW2	A	502	2	28,28,28	1.45	2 (7%)	37,39,39	1.68	8 (21%)
3	PW2	F	502	2	28,28,28	1.47	2 (7%)	37,39,39	1.61	7 (18%)
3	PW2	C	502	2	28,28,28	1.44	2 (7%)	37,39,39	1.46	6 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PW2	D	502	2	-	6/15/24/24	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PW2	E	502	2	-	4/15/24/24	0/3/3/3
3	PW2	B	502	2	-	4/15/24/24	0/3/3/3
4	GOL	E	503	-	-	0/4/4/4	-
3	PW2	A	502	2	-	4/15/24/24	0/3/3/3
3	PW2	F	502	2	-	6/15/24/24	0/3/3/3
3	PW2	C	502	2	-	6/15/24/24	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	502	PW2	C20-N21	5.16	1.45	1.33
3	A	502	PW2	C20-N21	5.09	1.45	1.33
3	F	502	PW2	C20-N21	5.07	1.44	1.33
3	B	502	PW2	C20-N21	4.97	1.44	1.33
3	D	502	PW2	C20-N21	4.88	1.44	1.33
3	C	502	PW2	C20-N21	4.86	1.44	1.33
3	B	502	PW2	O26-C20	-2.38	1.18	1.23
3	D	502	PW2	O26-C20	-2.38	1.18	1.23
3	F	502	PW2	O26-C20	-2.32	1.18	1.23
3	A	502	PW2	O26-C20	-2.24	1.18	1.23
3	C	502	PW2	O26-C20	-2.24	1.18	1.23
3	E	502	PW2	O26-C20	-2.14	1.18	1.23
3	B	502	PW2	O01-C02	2.03	1.40	1.36
3	E	502	PW2	O01-C02	2.01	1.40	1.36

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	502	PW2	C04-C03-C02	-5.47	116.32	120.68
3	A	502	PW2	C04-C03-C02	-4.75	116.90	120.68
3	B	502	PW2	C04-C03-C02	-4.32	117.24	120.68
3	B	502	PW2	C19-C20-N21	4.25	119.82	115.67
3	C	502	PW2	C04-C03-C02	-4.18	117.35	120.68
3	D	502	PW2	C19-C20-N21	4.18	119.74	115.67
3	F	502	PW2	C04-C03-C02	-4.13	117.39	120.68
3	F	502	PW2	C19-C20-N21	4.09	119.66	115.67
3	D	502	PW2	C03-C04-C17	3.93	121.01	117.11
3	E	502	PW2	C19-C20-N21	3.83	119.41	115.67
3	A	502	PW2	C03-C04-C17	3.79	120.87	117.11
3	E	502	PW2	C04-C03-C02	-3.55	117.86	120.68
3	B	502	PW2	C03-C04-C17	3.37	120.45	117.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	502	PW2	C06-C05-C04	-3.15	115.89	121.36
3	C	502	PW2	C19-C20-N21	3.13	118.72	115.67
3	E	502	PW2	C17-N18-C19	3.02	122.67	117.67
3	D	502	PW2	C17-C04-C05	-2.78	116.86	121.69
3	F	502	PW2	C03-C04-C17	2.75	119.83	117.11
3	A	502	PW2	C03-C04-C05	-2.68	116.43	120.86
3	A	502	PW2	C20-C19-N18	2.65	120.96	115.78
3	C	502	PW2	C03-C04-C17	2.59	119.68	117.11
3	C	502	PW2	C23-C22-N21	-2.59	105.08	113.06
3	A	502	PW2	C04-C17-N18	-2.58	120.07	124.32
3	F	502	PW2	C17-N18-C19	2.57	121.94	117.67
3	B	502	PW2	O11-C10-C09	-2.54	117.51	124.43
3	B	502	PW2	C04-C17-N18	-2.54	120.14	124.32
3	C	502	PW2	C17-N18-C19	2.52	121.86	117.67
3	B	502	PW2	C17-N18-C19	2.46	121.76	117.67
3	F	502	PW2	C06-C05-C04	-2.44	117.12	121.36
3	D	502	PW2	C04-C17-N18	-2.39	120.38	124.32
3	E	502	PW2	C04-C17-N18	-2.37	120.41	124.32
3	E	502	PW2	C03-C04-C17	2.37	119.45	117.11
3	F	502	PW2	C04-C17-N18	-2.34	120.45	124.32
3	D	502	PW2	C06-C05-C16	2.28	121.67	118.09
3	D	502	PW2	O11-C10-C09	-2.19	118.45	124.43
3	C	502	PW2	C04-C17-N18	-2.18	120.72	124.32
3	F	502	PW2	C23-C22-N21	-2.13	106.50	113.06
3	D	502	PW2	C12-O11-C10	-2.10	112.95	117.51
3	A	502	PW2	C17-N18-C19	2.10	121.15	117.67
3	A	502	PW2	O25-C23-C22	2.00	119.70	112.74

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	PW2	C02-C19-C20-N21
3	A	502	PW2	N18-C19-C20-N21
3	A	502	PW2	N18-C19-C20-O26
3	D	502	PW2	C09-C10-O11-C12
3	D	502	PW2	C13-C10-O11-C12
3	C	502	PW2	C09-C10-O11-C12
3	C	502	PW2	C13-C10-O11-C12
3	F	502	PW2	C13-C10-O11-C12
3	F	502	PW2	C09-C10-O11-C12
3	B	502	PW2	N18-C19-C20-O26

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Mol	Chain	Res	Type	Atoms
3	C	502	PW2	N18-C19-C20-O26
3	D	502	PW2	N18-C19-C20-O26
3	E	502	PW2	N18-C19-C20-O26
3	F	502	PW2	N18-C19-C20-O26
3	B	502	PW2	C02-C19-C20-N21
3	C	502	PW2	C02-C19-C20-N21
3	D	502	PW2	C02-C19-C20-N21
3	E	502	PW2	C02-C19-C20-N21
3	F	502	PW2	C02-C19-C20-N21
3	B	502	PW2	N18-C19-C20-N21
3	C	502	PW2	N18-C19-C20-N21
3	D	502	PW2	N18-C19-C20-N21
3	E	502	PW2	N18-C19-C20-N21
3	F	502	PW2	N18-C19-C20-N21
3	A	502	PW2	C02-C19-C20-O26
3	B	502	PW2	C02-C19-C20-O26
3	C	502	PW2	C02-C19-C20-O26
3	D	502	PW2	C02-C19-C20-O26
3	E	502	PW2	C02-C19-C20-O26
3	F	502	PW2	C02-C19-C20-O26

There are no ring outliers.

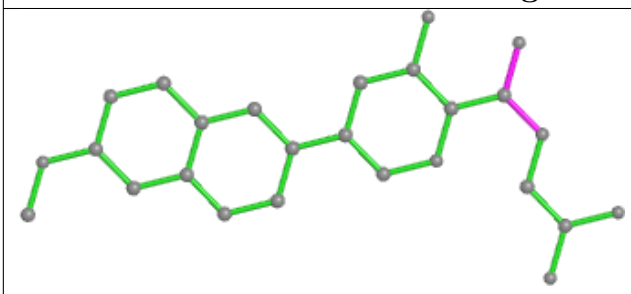
8 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	504	SO4	1	0
3	D	502	PW2	3	0
3	E	502	PW2	1	0
3	B	502	PW2	1	0
4	E	503	GOL	2	0
3	A	502	PW2	1	0
3	F	502	PW2	1	0
3	C	502	PW2	3	0

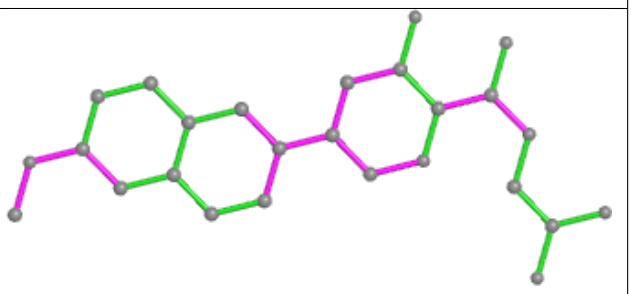
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

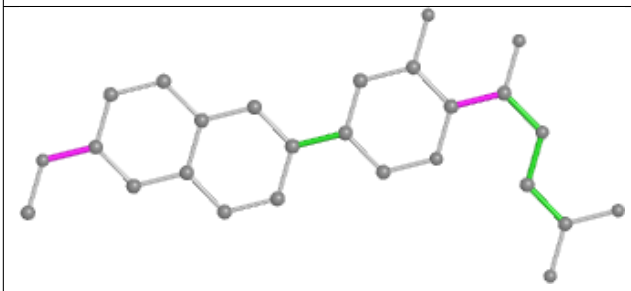
Ligand PW2 D 502



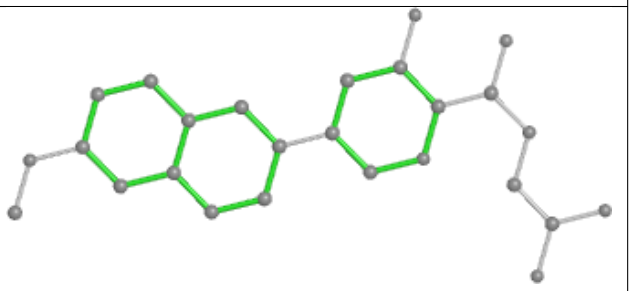
Bond lengths



Bond angles

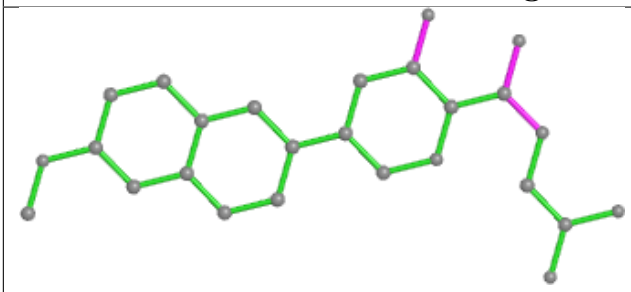


Torsions

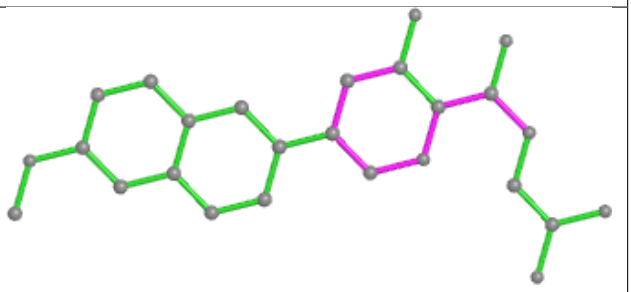


Rings

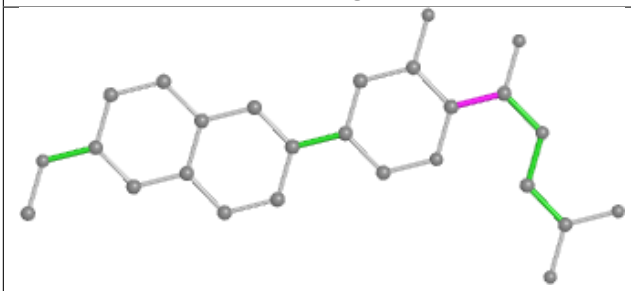
Ligand PW2 E 502



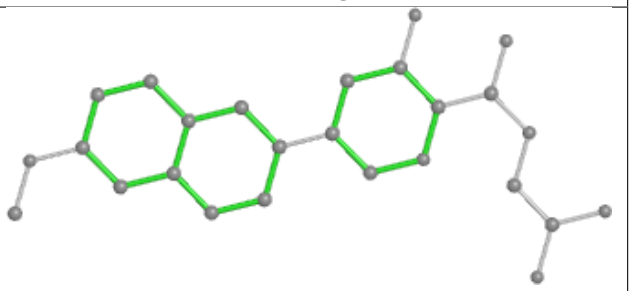
Bond lengths



Bond angles

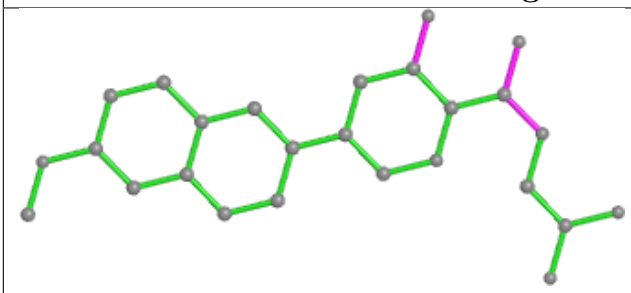


Torsions

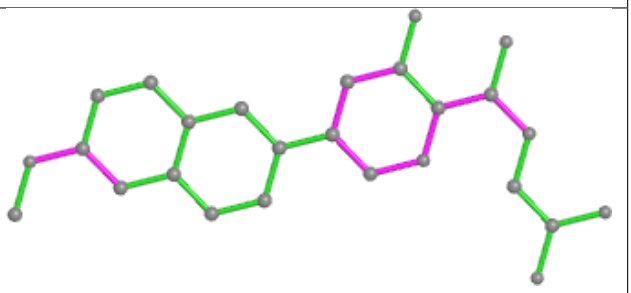


Rings

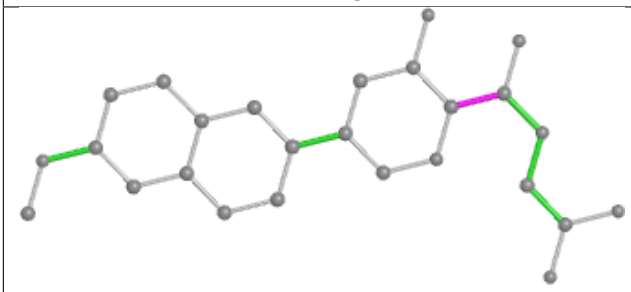
Ligand PW2 B 502



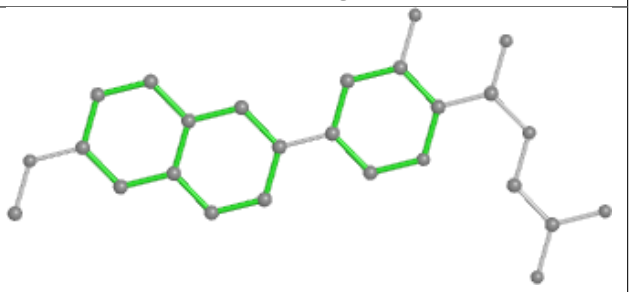
Bond lengths



Bond angles

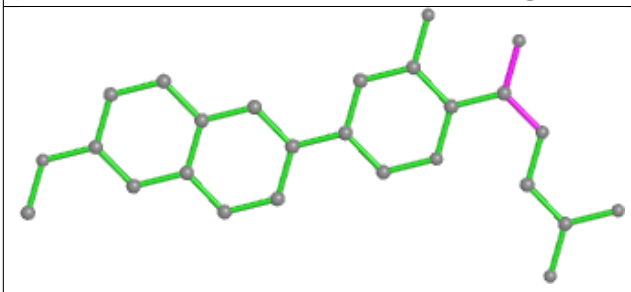


Torsions

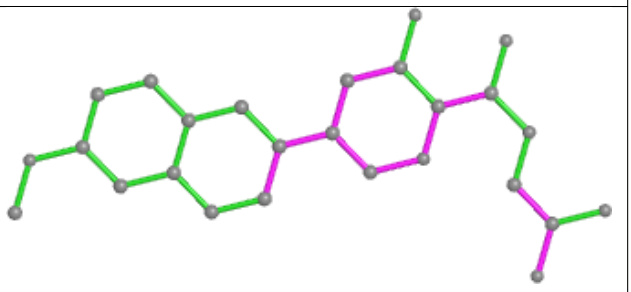


Rings

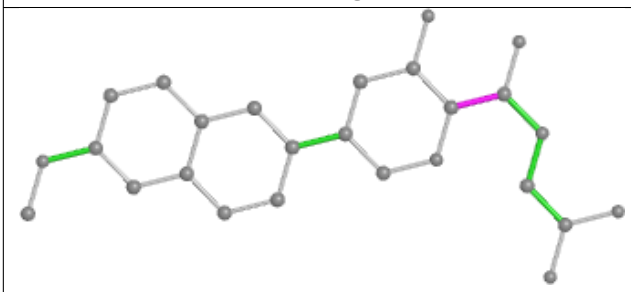
Ligand PW2 A 502



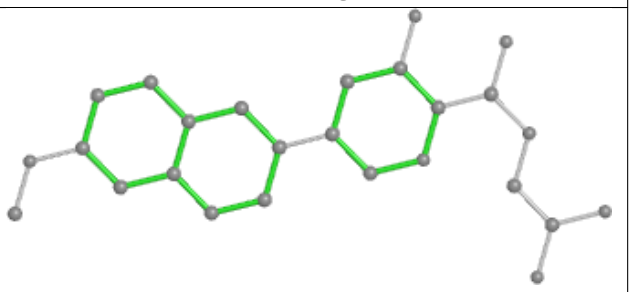
Bond lengths



Bond angles

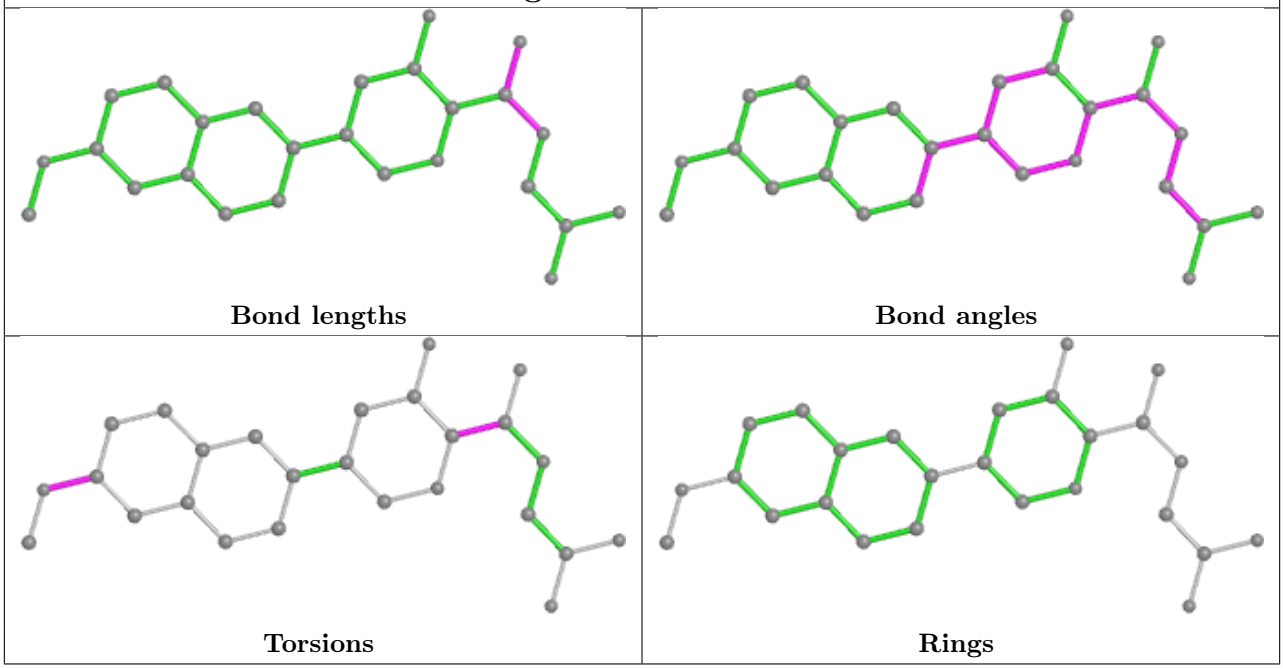


Torsions

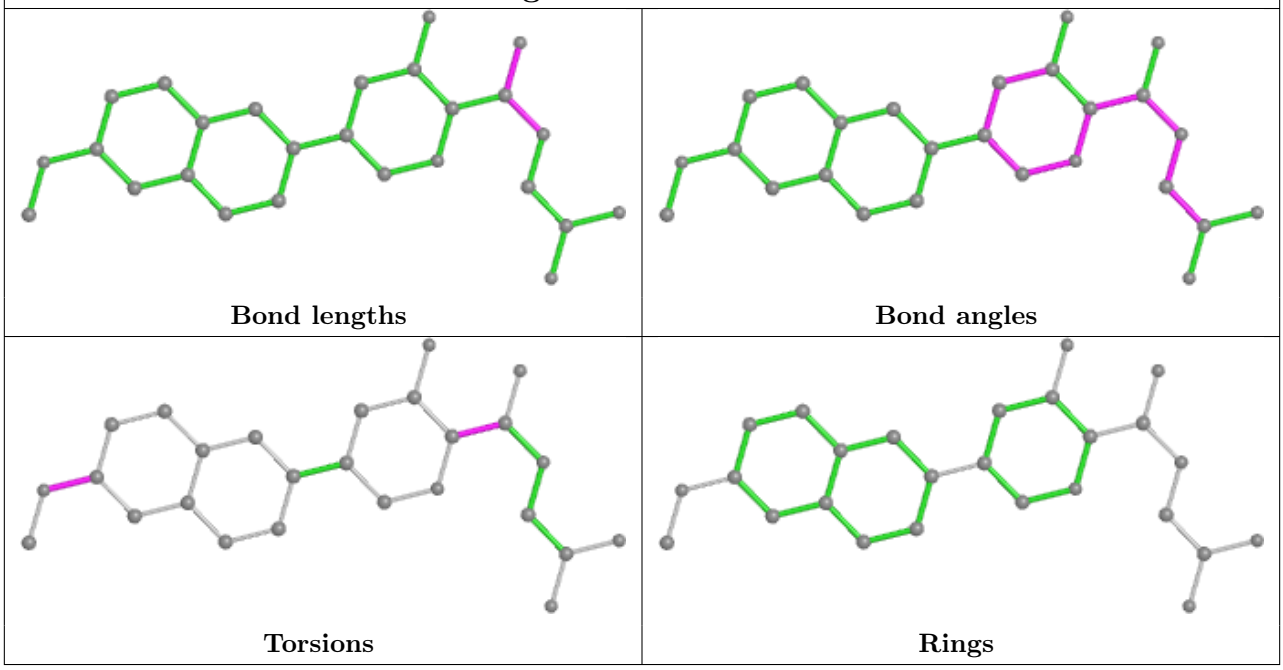


Rings

Ligand PW2 F 502



Ligand PW2 C 502



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	197/252 (78%)	0.36	13 (6%)	18	12	28, 49, 72, 91	0
1	B	199/252 (78%)	0.31	15 (7%)	14	9	31, 48, 75, 156	0
1	C	196/252 (77%)	0.39	13 (6%)	18	12	31, 47, 76, 122	0
1	D	199/252 (78%)	0.34	14 (7%)	16	11	30, 46, 71, 103	0
1	E	194/252 (76%)	0.33	12 (6%)	20	15	33, 48, 76, 129	0
1	F	197/252 (78%)	0.43	16 (8%)	12	7	31, 51, 79, 138	0
All	All	1182/1512 (78%)	0.36	83 (7%)	16	11	28, 48, 76, 156	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	250	ASP	5.4
1	B	350	LYS	5.2
1	E	349	GLY	4.9
1	C	326	CYS	4.8
1	C	349	GLY	4.7
1	B	351	ALA	4.7
1	D	351	ALA	4.6
1	A	254	ASP	4.5
1	F	351	ALA	4.5
1	E	351	ALA	4.4
1	D	326	CYS	4.2
1	B	349	GLY	4.1
1	D	352	GLN	4.1
1	A	351	ALA	3.9
1	B	348	GLU	3.8
1	F	230	HIS	3.7
1	C	351	ALA	3.7
1	E	326	CYS	3.5
1	A	333	ASP	3.5

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Mol	Chain	Res	Type	RSRZ
1	F	352	GLN	3.5
1	A	230	HIS	3.2
1	F	231	ASP	3.1
1	C	325	THR	3.1
1	A	326	CYS	3.1
1	B	253	GLY	3.1
1	E	304	PRO	3.1
1	C	254	ASP	3.0
1	C	387	THR	3.0
1	F	333	ASP	3.0
1	B	335	ASP	2.9
1	B	403	TYR	2.9
1	E	352	GLN	2.8
1	E	350	LYS	2.8
1	D	232	THR	2.7
1	F	232	THR	2.7
1	E	325	THR	2.7
1	F	326	CYS	2.7
1	A	261	GLY	2.7
1	B	251	ILE	2.7
1	A	220	GLN	2.7
1	B	326	CYS	2.7
1	C	350	LYS	2.6
1	F	353	PHE	2.6
1	E	403	TYR	2.6
1	D	387	THR	2.6
1	D	231	ASP	2.6
1	E	284	ASN	2.5
1	F	262	LYS	2.5
1	A	257	THR	2.5
1	F	285	GLY	2.5
1	E	232	THR	2.5
1	D	403	TYR	2.5
1	F	349	GLY	2.5
1	B	306	ASN	2.5
1	A	231	ASP	2.5
1	D	251	ILE	2.4
1	A	306	ASN	2.4
1	D	327	ILE	2.4
1	B	325	THR	2.4
1	F	350	LYS	2.4
1	C	284	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	327	ILE	2.3
1	A	232	THR	2.3
1	C	388	VAL	2.3
1	F	387	THR	2.3
1	A	387	THR	2.3
1	D	230	HIS	2.2
1	C	353	PHE	2.2
1	B	220	GLN	2.2
1	B	254	ASP	2.2
1	E	327	ILE	2.2
1	C	260	GLU	2.2
1	C	289	SER	2.2
1	E	348	GLU	2.2
1	F	197	TYR	2.1
1	B	231	ASP	2.1
1	D	349	GLY	2.1
1	F	289	SER	2.1
1	A	188	LEU	2.1
1	D	335	ASP	2.1
1	F	306	ASN	2.1
1	D	353	PHE	2.1
1	B	188	LEU	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
3	PW2	A	502	26/26	0.82	0.32	31,47,70,74	0

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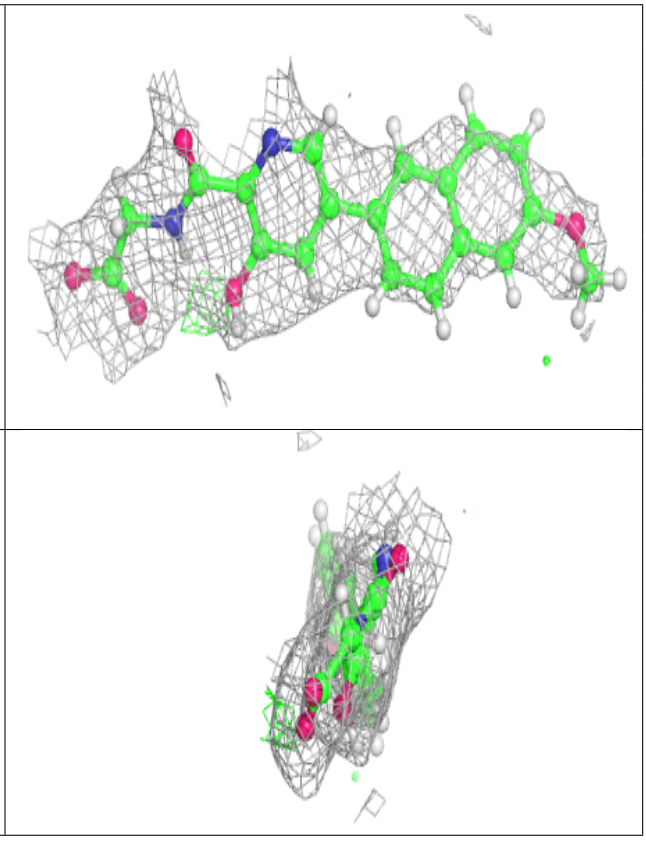
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	PW2	F	502	26/26	0.82	0.29	34,42,65,67	0
3	PW2	C	502	26/26	0.83	0.30	31,41,56,57	0
3	PW2	E	502	26/26	0.85	0.27	27,37,51,56	0
3	PW2	B	502	26/26	0.88	0.27	32,46,61,67	0
4	GOL	E	503	6/6	0.88	0.21	41,50,60,63	0
3	PW2	D	502	26/26	0.89	0.25	31,39,65,68	0
5	SO4	E	504	5/5	0.89	0.35	66,71,111,117	0
2	MN	A	501	1/1	0.98	0.21	36,36,36,36	0
2	MN	E	501	1/1	0.99	0.15	31,31,31,31	0
2	MN	F	501	1/1	0.99	0.16	36,36,36,36	0
2	MN	B	501	1/1	0.99	0.17	29,29,29,29	0
2	MN	C	501	1/1	0.99	0.20	32,32,32,32	0
2	MN	D	501	1/1	0.99	0.16	26,26,26,26	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.

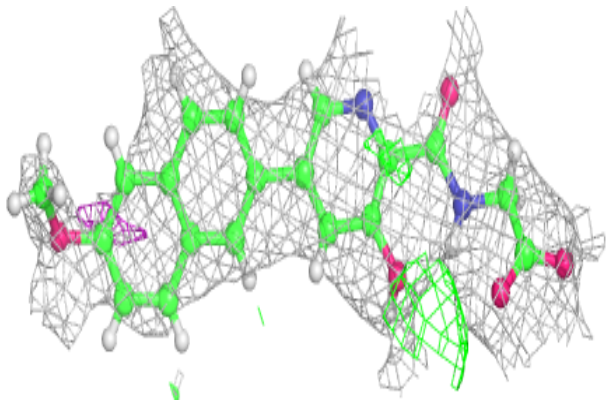
Electron density around PW2 A 502:

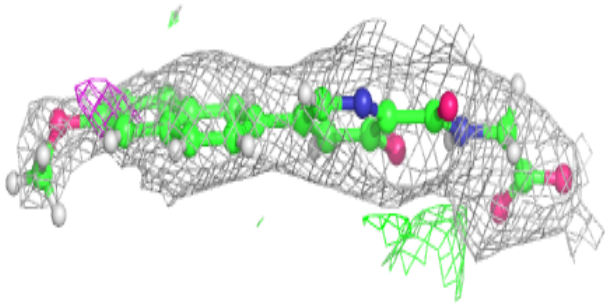
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

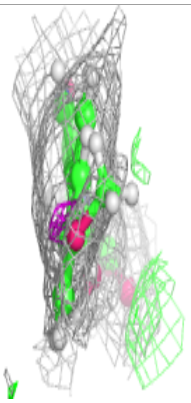


Electron density around PW2 F 502:

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

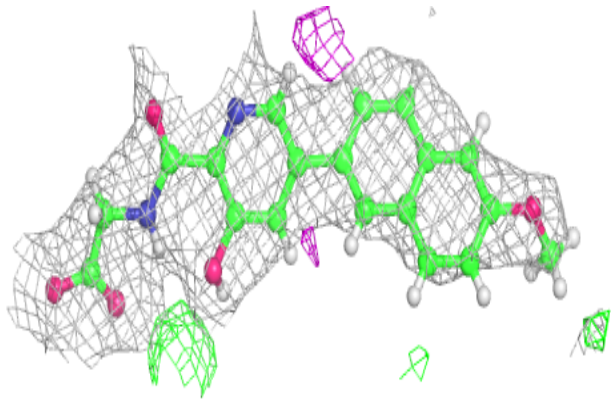


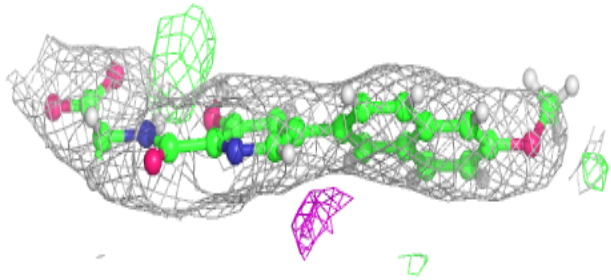


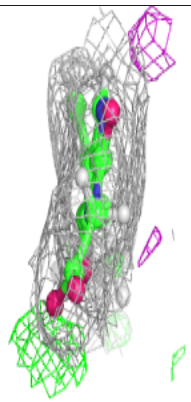


Electron density around PW2 C 502:

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

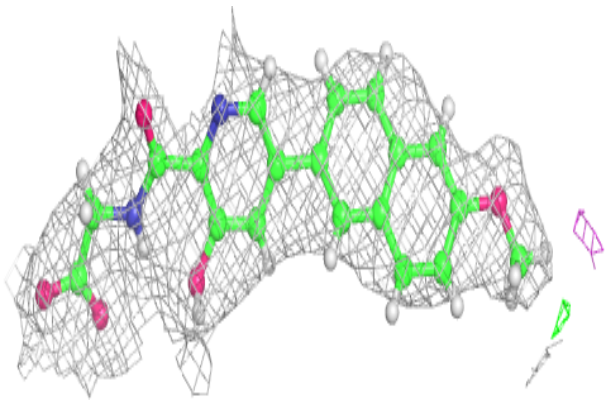


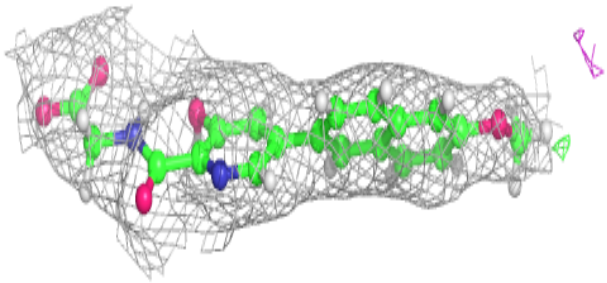


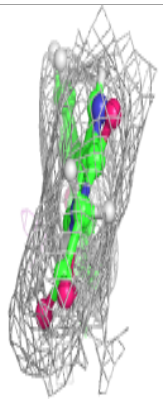


Electron density around PW2 E 502:

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

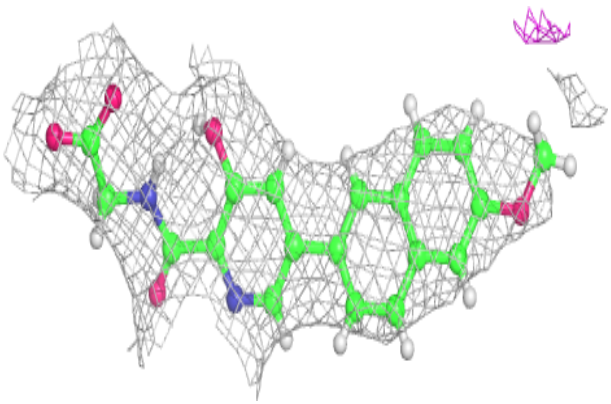


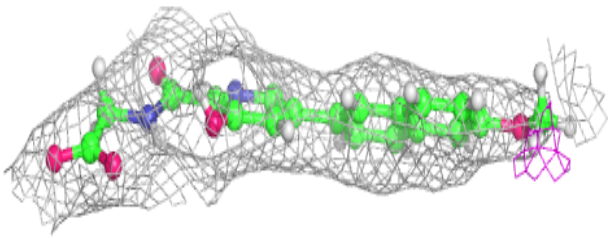


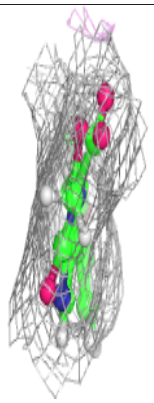


Electron density around PW2 B 502:

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

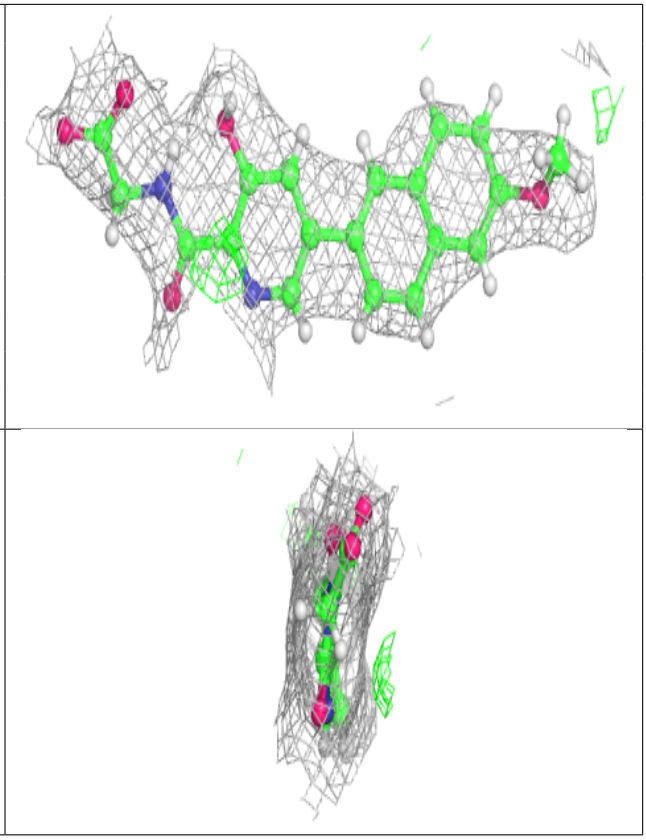






Electron density around PW2 D 502:

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.