

AHEAD-2020 WP14

Tasks 14.6 & 14.8

- Blind-Line-Search and ID (14.6)
- Time-Evolving Photoionization (14.8)

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Blind Line Search Routine

- Fortran-90 Routine working in either HEASOFT or CIAO environments
- Extremely fast on 30,000 channel spectra (e.g. Athena-XIFU)
- Reads in:
 - RMF/RSP (FITS format)
 - Continuum-Normalized data in sigma (QDP|ASCII: e.g. from XSPEC or Sherpa)
 - Output Filenames prefix
 - (line statistical significance threshold in sigma + verbosity: default $\sigma=4$, $v=0$)
- Extracts LSF, computes LSF FWHM (i.e. spectral resolution)
- Scans the data searching for positive/negative unresolved, quasi-symmetric, line-like features, with integrated significance exceeding the threshold
- Reads in atomic-database (currently Verner et al., 2006) and looks for possible line-IDs (WiP)
- Writes out:
 - File containing list of 'detected' lines' energy, peak-significance, integrated-significance
 - File containing possible redshift IDs (WiP)

```
nicastro@Falcao.local% bls_v2 athena_xifu_1469_onaxis_pitch249um_v20160401.rsp LSF-convolved_Delchi_28_1.5e7.dat 4 0
```

```
Response Filename: athena_xifu_1469_onaxis_pitch249um_v20160401.rsp
Continuum-Normalized Filename: LSF-convolved_Delchi_28_1.5e7.dat
Output Filename: 4
Line Threshold Significance: 0
Using default Verbosity: 0
```

```
Number of Absorption Lines:      18
Number of Emission Lines:       2
```

BLS Example: "Observation"

Line of Sight Extracted from Cen & Ostriker (2006) Hydro-Dynamical Simulations

Column Density

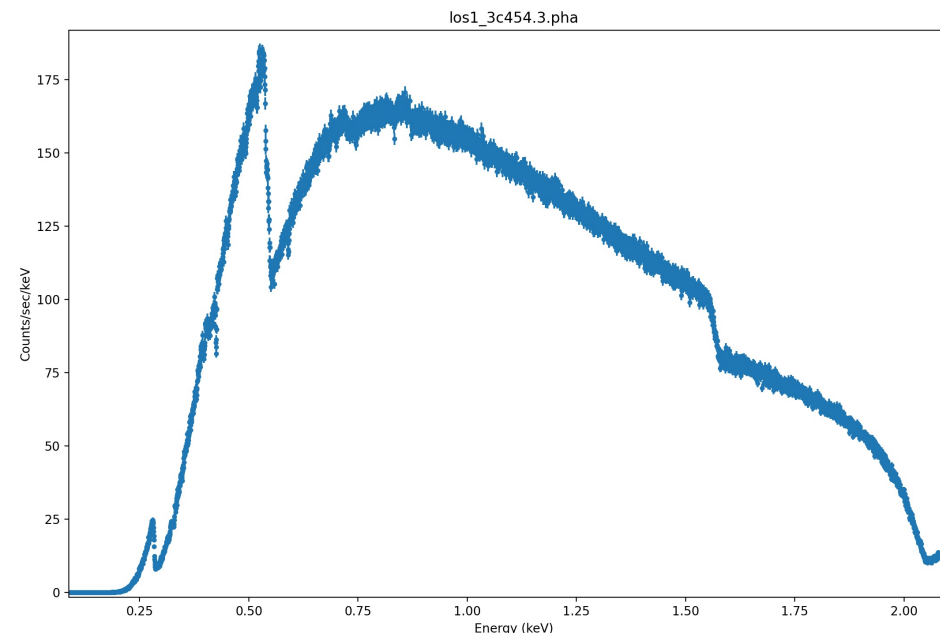
Redshift

Metallicity

```
nicastro@Falcao.local% more los_28_WoZ_WHIM.out
```

No. of Filament, Last row of the Group in Original File; No of rows in Group; T(in K); NH(in cm-2); Turbulence (in km/s); Redshift; log(U); Density (at the given z); Metallicity (compared to Solar)									
1	3847	1	1379361.25	3.36197258E+18	261.441772	0.109475948	-1.93419802	1.08900003E-05	9.50078666E-02
2	4908	1	729345.250	3.23784322E+18	190.108902	0.140729979	-1.91785955	1.04900000E-05	0.201572582
3	4915	3	688116.562	3.43952801E+18	184.602371	0.140762627	-1.94454646	1.11622676E-05	0.506648600
4	4919	2	907217.188	3.20123498E+18	212.001709	0.140754178	-1.92046666	1.05629624E-05	0.220032334
5	9167	3	206526.281	3.33116976E+18	101.091904	0.273898929	-1.94133806	1.10789997E-05	1.66198611E-02
6	13305	6	214686.875	1.11909778E+19	103.111465	0.416768640	-2.30846786	3.37445308E-05	2.01087464E-02
7	13919	25	5417012.50	1.18929300E+20	515.280945	0.436983317	-2.99046493	7.81381386E-04	0.115085848
8	15462	1	115829.562	1.46608968E+19	75.7609711	0.495346159	-2.57376432	4.75099987E-05	1.01446602E-02
9	17371	7	451797.094	5.50937954E+18	149.567551	0.569676399	-2.06023479	1.45867716E-05	0.191525415
10	17940	4	143990.109	4.48879305E+18	84.4650497	0.590543449	-2.10365844	1.60944182E-05	1.46267880E-02
11	18372	4	398528.188	4.00824627E+18	140.527100	0.608431876	-2.02583265	1.34562370E-05	2.12974884E-02
12	18414	6	286541.844	1.97825693E+19	119.021568	0.609701395	-1.99990737	1.45019585E-05	0.195985019
13	18419	4	371265.344	1.66751582E+19	135.493530	0.610535920	-2.91440797	1.06934771E-04	0.309740871
14	19595	3	1126450.88	3.17020483E+18	236.260986	0.659921110	-1.91196179	1.03496823E-05	0.154758915
15	21110	5	148218.734	3.50232332E+18	85.6894989	0.724813461	-1.91181695	1.03493585E-05	7.48509616E-02

Athena-XIFU 100 ks mock spectrum of 3C454.3 (z=0.859)
@1 mCrab



BLS Example: output

Tentative Redshift IDs

Detected Line-list

```
# Number of Absorption Lines:      17
# Number of Emission Lines:        2
# E (keV), Sigma-Peak, Sigma, Sigma-Threshold
# Negative Lines:
0.267800003    1.59371269    4.03757238    4.00000000
0.356200010    7.09524679    17.2754307    4.00000000
0.413399994    1.77576268    4.03026819    4.00000000
0.454200000    10.0198488    24.6709194    4.00000000
0.503000021    8.24497032    20.4788818    4.00000000
0.564999998    2.81150985    7.20422220    4.00000000
0.573800027    15.4453526    39.6384010    4.00000000
0.606999993    3.29755354    8.53451443    4.00000000
0.623799980    1.95347917    4.41314697    4.00000000
0.640600026    5.42002916    7.90693521    4.00000000
0.641399980    5.75636578    14.7684107    4.00000000
0.710600019    7.27873373    18.1650715    4.00000000
0.711399972    6.76571894    9.38497066    4.00000000
0.783399999    1.63223302    4.25715160    4.00000000
0.940599978    6.63401318    17.7794666    4.00000000
 1.29740000    5.49634790    15.2450886    4.00000000
 1.71099997    1.81207728    4.86538553    4.00000000
# Positive Lines:
0.470999986    1.57258713    4.53434849    4.00000000
0.685000002    1.54889691    4.23500299    4.00000000
```

```
Possible Redshift: z = 0.611312509
=====
1: Eobs = 0.356200010 ; E0 = 0.573949516 ; Ion: O7
 2: Eobs = 0.413399994 ; E0 = 0.666116595 ; Ion: O7
 3: Eobs = 0.573800027 ; E0 = 0.924571157 ; Ion: Ne9
 4: Eobs = 0.623799980 ; E0 = 1.00513673 ; Ion: Fe21
```

```
Possible Redshift: z = 0.388363600
=====
1: Eobs = 0.413399994 ; E0 = 0.573949516 ; Ion: O7
 2: Eobs = 0.503000021 ; E0 = 0.698346913 ; Ion: O7
 3: Eobs = 0.623799980 ; E0 = 0.866061211 ; Ion: Fe18
 5: Eobs = 0.640600026 ; E0 = 0.889385760 ; Ion: Fe18
 6: Eobs = 0.641399980 ; E0 = 0.890496373 ; Ion: Fe18
```

```
Possible Redshift: z = 0.263649344
=====
1: Eobs = 0.454200000 ; E0 = 0.573949516 ; Ion: O7
 2: Eobs = 0.564999998 ; E0 = 0.713961899 ; Ion: O7
 3: Eobs = 0.710600019 ; E0 = 0.897949219 ; Ion: Fe17
 4: Eobs = 0.711399972 ; E0 = 0.898960114 ; Ion: Fe17
```

```
Possible Redshift: z = 0.141052723
=====
1: Eobs = 0.503000021 ; E0 = 0.573949516 ; Ion: O7
 2: Eobs = 0.573800027 ; E0 = 0.654736102 ; Ion: O8
 4: Eobs = 0.623799980 ; E0 = 0.711788654 ; Ion: O7
 5: Eobs = 0.783399999 ; E0 = 0.893900692 ; Ion: Fe17
 6: Eobs = 0.940599978 ; E0 = 1.07327414 ; Ion: Ne9
```

```
Possible Redshift: z = 1.58398151E-02
=====
1: Eobs = 0.564999998 ; E0 = 0.573949516 ; Ion: O7
 2: Eobs = 0.640600026 ; E0 = 0.650747001 ; Ion: O8
 4: Eobs = 0.641399980 ; E0 = 0.651559651 ; Ion: O8
 6: Eobs = 0.710600019 ; E0 = 0.721855819 ; Ion: O7
 7: Eobs = 0.711399972 ; E0 = 0.722668409 ; Ion: O7
```

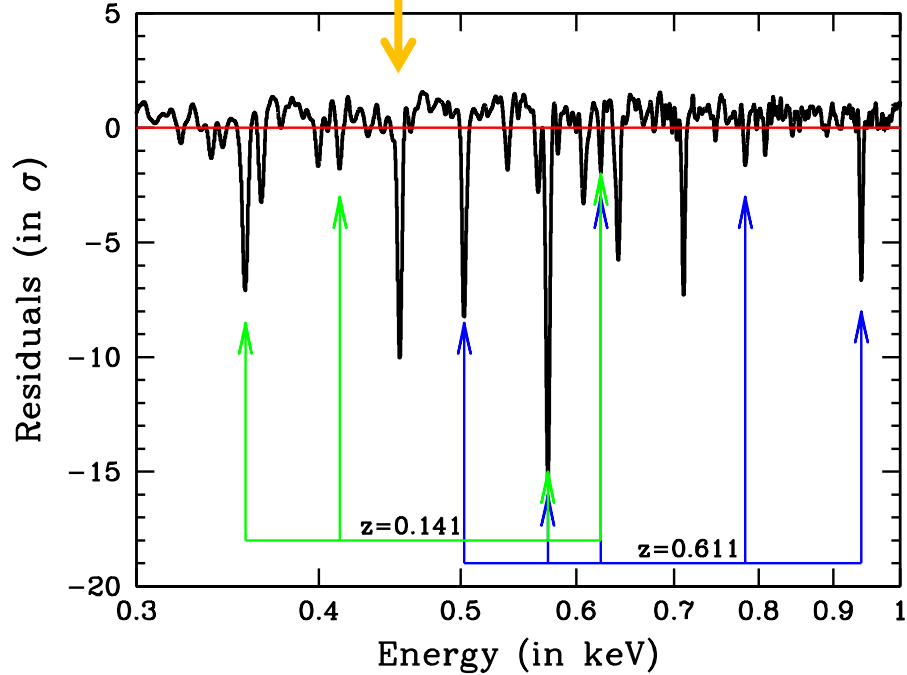
```
Possible Redshift: z = 2.60472298E-04
=====
1: Eobs = 0.573800027 ; E0 = 0.573949516 ; Ion: O7
 2: Eobs = 0.710600019 ; E0 = 0.710785091 ; Ion: O7
 3: Eobs = 0.711399972 ; E0 = 0.711585283 ; Ion: O7
```

```
Possible Redshift: z = -5.44489026E-02
=====
1: Eobs = 0.606999993 ; E0 = 0.573949516 ; Ion: O7
 2: Eobs = 0.940599978 ; E0 = 0.889385343 ; Ion: Fe18
```

```
Possible Redshift: z = -7.99141526E-02
=====
1: Eobs = 0.623799980 ; E0 = 0.573949516 ; Ion: O7
 2: Eobs = 0.710600019 ; E0 = 0.653813004 ; Ion: O8
 4: Eobs = 0.711399972 ; E0 = 0.654549062 ; Ion: O8
 6: Eobs = 0.783399999 ; E0 = 0.720795274 ; Ion: O7
 7: Eobs = 0.940599978 ; E0 = 0.865432739 ; Ion: Fe18
```

BLS Example: Line and Redshift IDs

Missed ID: OVIII Ly α z=0.437



Possible Redshift: $z = 0.611312509$

```
=====
1: Eobs = 0.356200010 ; E0 = 0.573949516 ; lon: O7
2 : Eobs = 0.413399994 ; E0 = 0.666116595 ; lon: O7
3 : Eobs = 0.573800027 ; E0 = 0.924571157 ; lon: Ne9
4 : Eobs = 0.623799980 ; E0 = 1.00513673 ; lon: Fe21
```

Possible Redshift: $z = 0.141052723$

```
=====
1: Eobs = 0.503000021 ; E0 = 0.573949516 ; lon: O7
2 : Eobs = 0.573800027 ; E0 = 0.654736102 ; lon: O8
4 : Eobs = 0.623799980 ; E0 = 0.711788654 ; lon: O7
5 : Eobs = 0.783399999 ; E0 = 0.893900692 ; lon: Fe17
6 : Eobs = 0.940599978 ; E0 = 1.07327414 ; lon: Ne9
```

BLS: To do (depending on funding, personnel, etc.)

- Update/Integrate (forbidden/intercombination, innershell resonant ground-state and metastable transitions) Atomic Database
- Improve Line/Redshift ID routine: needs to be smarter (AI training might help): e.g. starting from strongest and looking for right strength ratios (including saturation)
- Python interface to directly call BLS from within Sherpa | XSPEC
- GUI interface to locally iterate the procedure and plot results

14.8: Time-Evolving Photo-Ionization Device (TEPID: Luminari+22)

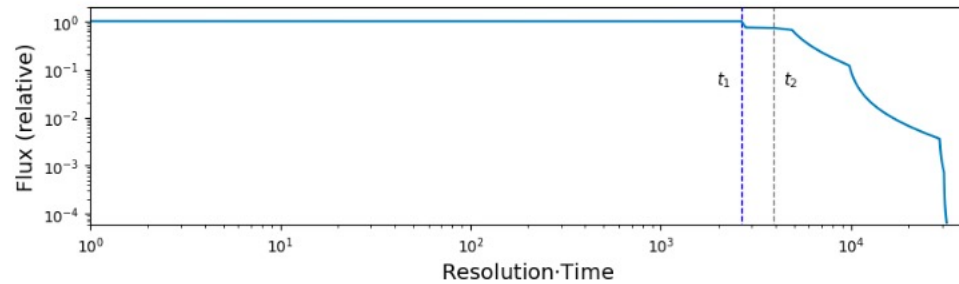
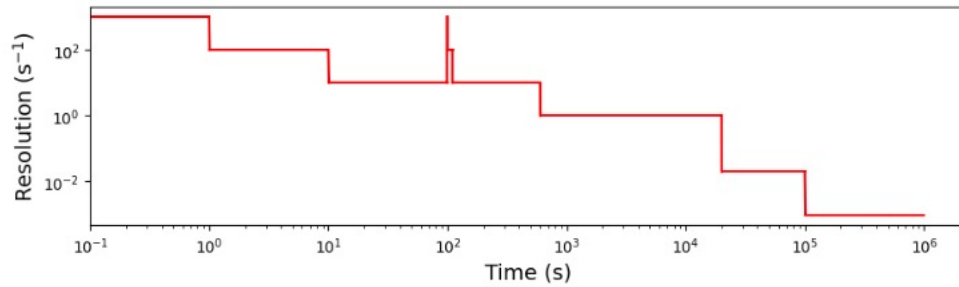
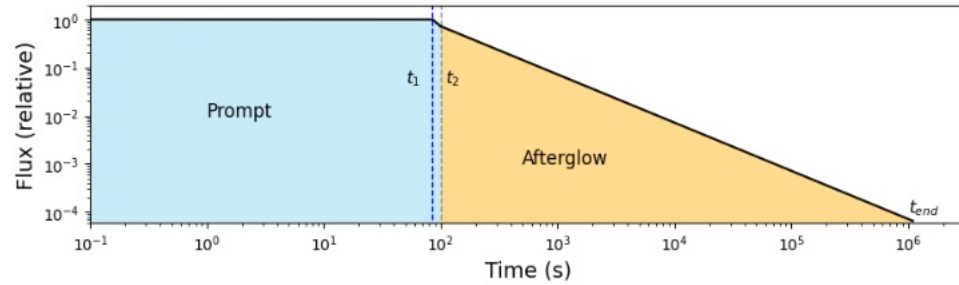
- Solves the system of 1st-order differential equations of time-evolving ionization balances following ionizing intensity variations with time
- Uses adaptive time-resolution algorithm to speed-up calculation
- Currently considers only photo-ionization and radiative recombinations and approximates on heating-cooling balance
- Includes radiative-transfer
- Applies to all variable ionizing sources: e.g. transients (GRBs) and AGNs
- Interfaces with Phase (Krongold+03), in XSPEC|Sherpa to produce mock spectra and fit data

$$\frac{dn_{X^i}}{dt} = -[F_{X^i} + \alpha_{\text{rec}}(X^{i-1}, T_e)n_e]n_{X^i} + F_{X^{i-1}}n_{X^{i-1}} + \alpha_{\text{rec}}(X^i, T_e)n_en_{X^{i+1}} \cdot \quad (4)$$

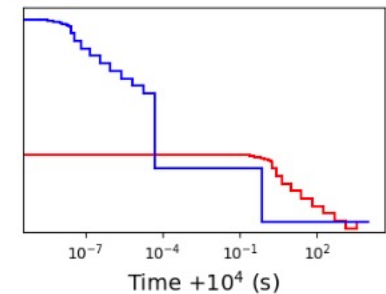
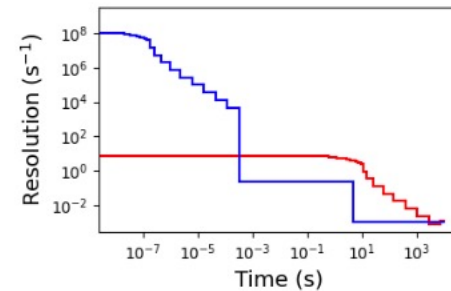
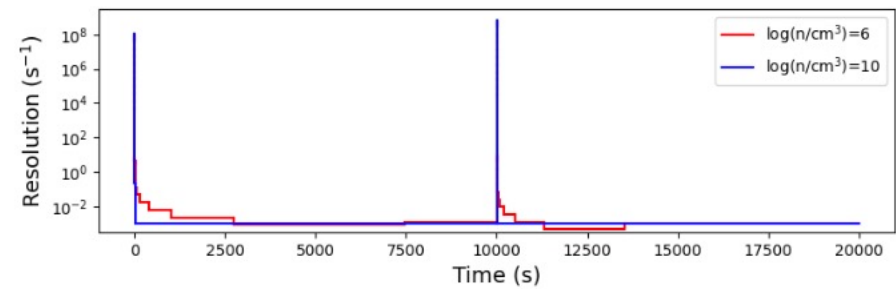
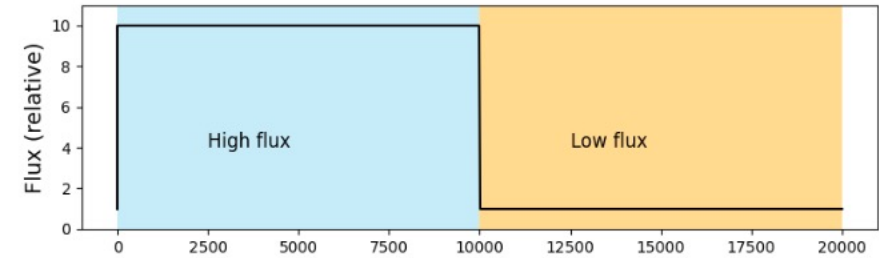
$$t_{eq} \approx \left\{ (\alpha_{rr}(X^i, T_e)n_e) \cdot \left(\frac{\alpha_{rr}(X^{i-1}, T_e)}{\alpha_{rr}(X^i, T_e)} + \frac{n_{X^{i+1}}}{n_{X^i}} \right) \right\}^{-1}$$

TEPID: Adaptive-Resolution Algorithm

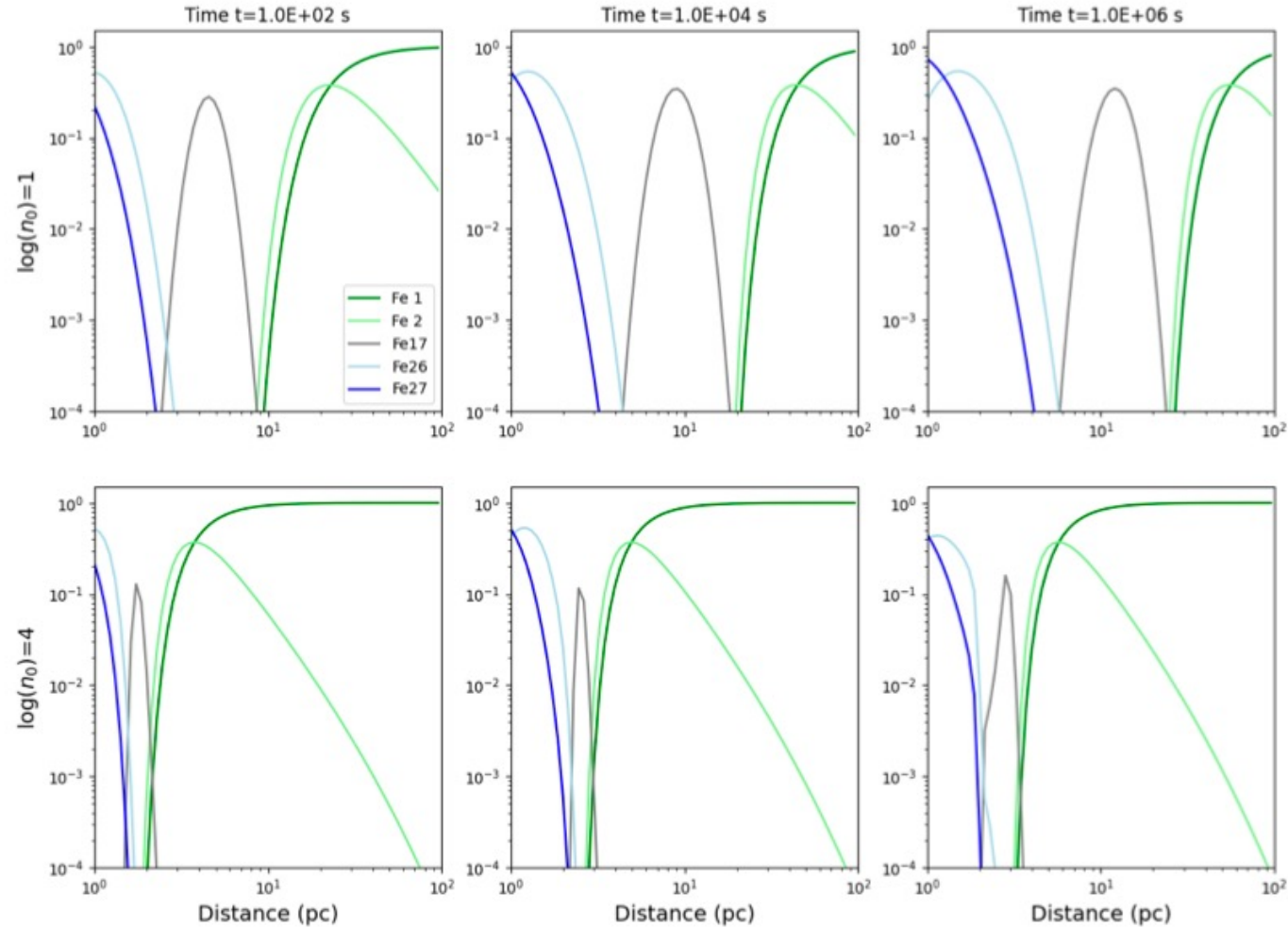
GRB case



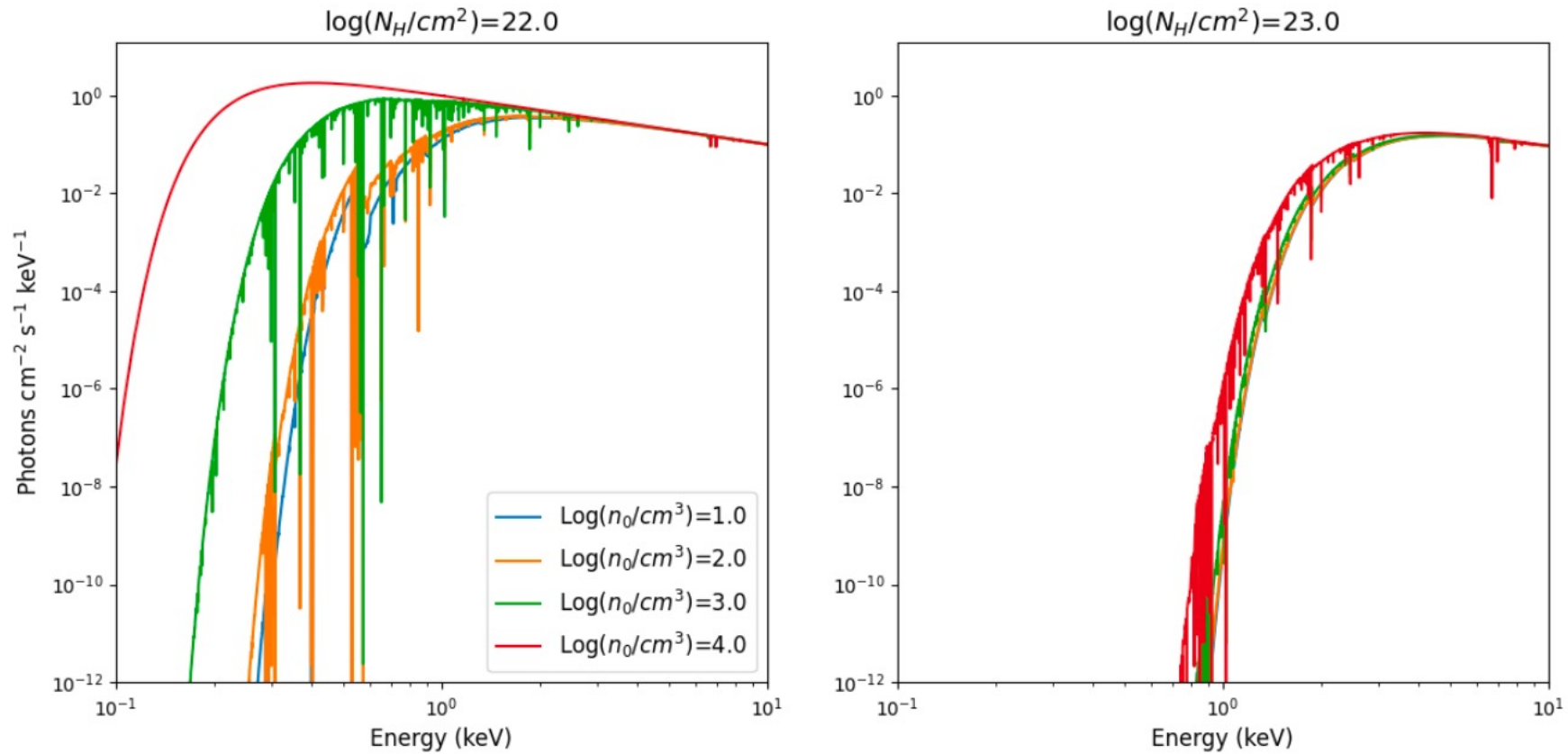
Up-and-Down case



TEPID: GRB-case: Stratified-Ionization



TEPID: GRB-case: Spectra vs Local Density



TEPID: AGN-case: Ionization Balance vs Time

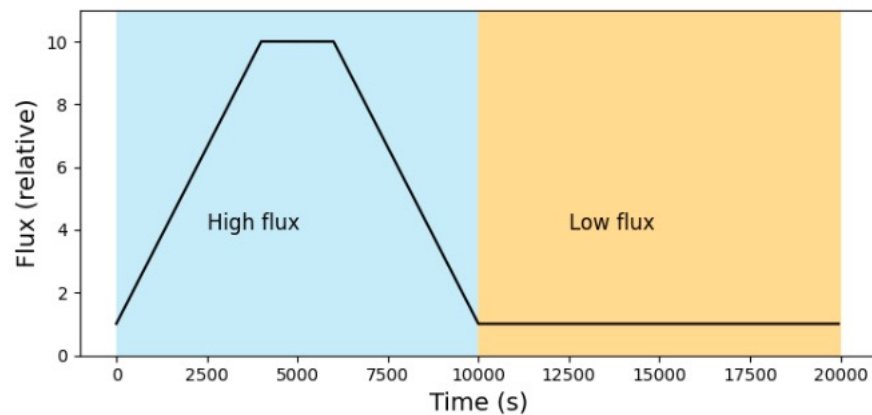
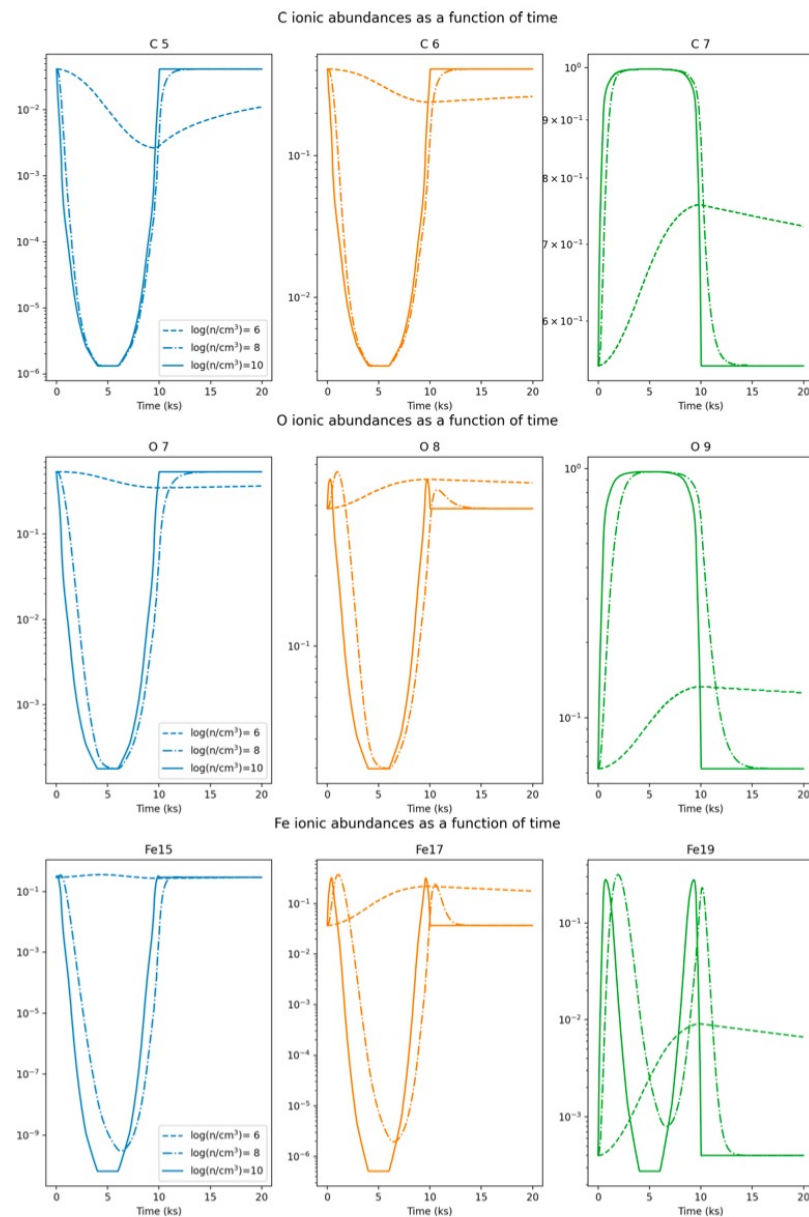
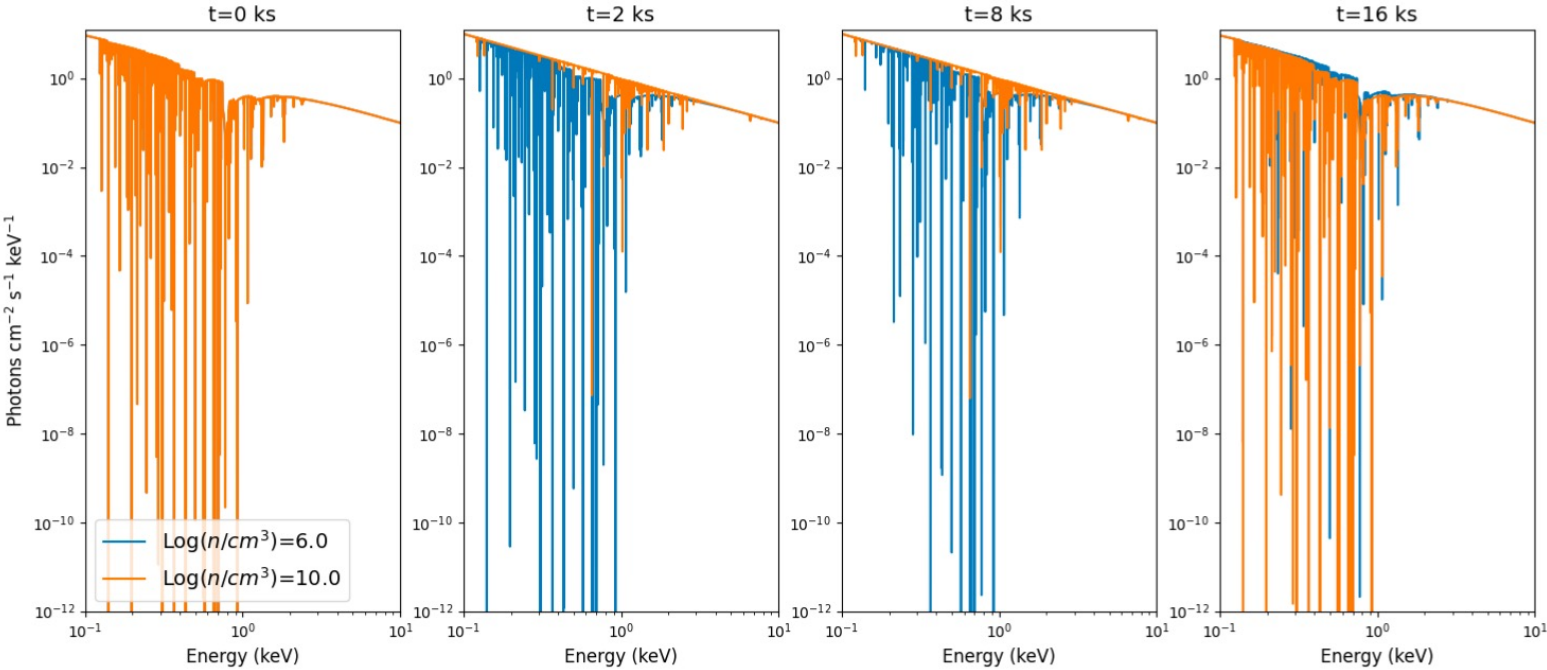
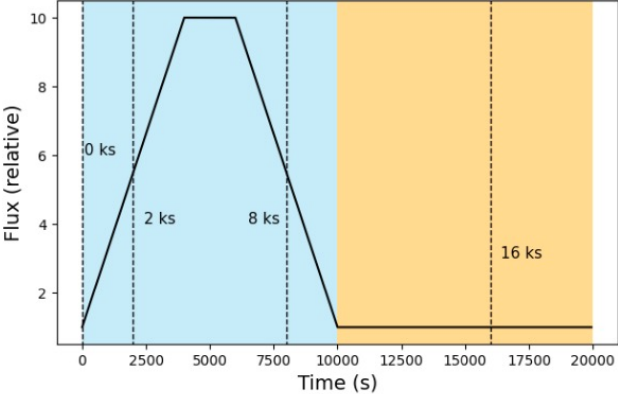


Fig. 9: AGN lightcurve



TEPID: AGN-case: Time-Resolved Spectra



TEPID: To do

- Include time-evolving population-level calculation and metastable transitions (to further constrain density)
- Include proper heating/cooling calculation
- Include additional ionization/recombination mechanisms (i.e. transition rates)
- Improve atomic-database
- Include proper pre-burst star-forming region ionization conditions based on simulations (Luminari, Graziani, in prep.)